

# **Transmittal**

**DATE:** February 12, 2009

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VIA: USMail

# WE ARE SENDING THE FOLLOWING MATERIALS:

Data validation report for Quendall Terminal

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# **DATA VALIDATION REPORT**

# Port of Seattle BNSF ROW - Quendall Terminal

# Prepared for:

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# Prepared by:

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February 9, 2009

Approved for Release:

Chris Ransom
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# **PROJECT NARRATIVE**

#### Basis for the Data Validation

This report summarizes the results of the full (Level IV) validation performed on water, soil, and sediment samples and the associated field and laboratory quality control samples collected for the Port of Seattle BNSF Right of Way project.

The samples were analyzed by Analytical Resources, Inc., Tukwila, Washington. The analytical methods and EcoChem project chemists are listed in the table below.

# ANALYSIS METHODS AND ECOCHEM CHEMISTS

Analysis	Method	Primary Review	Secondary Review
Semivolatile Organic Compounds (SVOC)	EPA SW8270D	Jennifer Newkirk	Christine Ransom
Polynuclear Aromatic Hydrocarbons (PAH)	EPA SW8270D-SIM	Jennifer Newkirk	Christine Ransom
PCB Aroclors	EPA SW8082	Melissa Swanson	Christine Ransom
Pesticides	EPA SW8081A	Melissa Swanson	Christine Ransom
Herbicides	EPA SW8151A	Melissa Swanson	Christine Ransom
Diesel Range Hydrocarbons	NWTPH-Dx	Linda Holz	Christine Ransom
Gasoline Range Hydrocarbons and BTEX Compounds	NWTPH-Gx, SW8121B	Linda Holz	Christine Ransom
Metals	EPA SW6020, SW7471A	Linda Holz	Christine Ransom
Conventionals (turbidity, percent solids)	EPA 180.1, 160.3	Linda Holz	Christine Ransom

The data were reviewed using guidance and quality control criteria documented in the analytical methods; the project quality assurance project plan (QAPP) *Port of Seattle, Rail Acquisition Study – Quendall Segment Right of Way* (Aug 8, 2008), and *National Functional Guidelines for Organic Data Review* (USEPA 1999) *and Inorganic Data Review* (1994,2004).

Following the **Summary of Findings**, a **Sample Index** and detailed validation reports are grouped by Transect or Boring event. Data Qualifier Definitions and Criteria Tables are included in **Appendix A**. **Appendix B** contains the Qualified Data Summary Table. **Appendix C** contains communication records. Data validation worksheets are kept on file at EcoChem.

#### **SUMMARY OF FINDINGS**

Overall, the data were acceptable. Out of a total of 4955 analytical results, one result for dinoseb was rejected, 149 results were estimated (J/UJ), and five results were qualified as not-detected (U) at an elevated reporting limit. Completeness for this data set is 99.8%. More than one set of results were reported for several samples due to re-extractions and necessary dilutions. In order to only have one result per analyte for each sample, data that should not be used were labeled do-not-report (DNR). Completeness is unaffected by the assigning of DNR flags.

Data quality issues that resulted in the qualification of data are noted below.

# **Baxter Mill Transect B1**

This transect consisted of 11 soil samples. One set of field replicates, B1-A-1.0 and DS-1, was submitted.

**SVOC:** The reporting limits for benzyl alcohol were estimated (UJ) in seven samples due to a continuing calibration (CCAL) percent difference (%D) outlier which indicated a low bias in the results.

**NWTPH-Dx:** The diesel range organic results for samples B1-A-1.0 and DS-1 were estimated (J) based on a field precision outlier.

*Metals:* There were no quality control issues that resulted in qualification of data.

# **Quendall Terminal North Transect Q1**

This transect consisted of one surface water, one sediment, and 13 soils. One set of field replicates, Q1-D-23.0 and SD-2, was submitted.

**SVOC:** The benzoic acid result for Sample Q1-D-3.5 was estimated (UJ) due to a low matrix spike (MS) recovery. The results for eight analytes (benzo(a)anthracene, benzo(b)fluoranthene, chrysene, indeno(1,2,3-cd)pyrene, benzo(a)pyrene, benzo(g,h,i)perylene, fluoranthene, and pyrene) in Samples Q1-D-23.0 and SD-2 were estimated (J) based on field precision outliers.

**Pesticides:** There were no quality control issues that resulted in qualification of data.

**PCB:** There were no quality control issues that resulted in qualification of data.

*Herbicides:* The dinoseb result for Sample Q1-Sed was rejected based on laboratory control sample and MS recoveries less than 10%.

**NWTPH-Dx:** The diesel range organic and motor oil results for Samples Q1-D-23.0 and SD-2 were estimated (J) based on field precision outliers.

**NWTPH-G:** There were no quality control issues that resulted in qualification of data.

**BTEX:** There were no quality control issues that resulted in qualification of data.

*Metals:* There were no quality control issues that resulted in qualification of data.

**Conventionals:** There were no quality control issues that resulted in qualification of data.

#### Quendall Terminal South Transect Q2 Quendall Terminal South

This transect consisted of one groundwater and 14 soils. One set of field replicates, Q2-D-W and WD-1, was submitted.

**SVOC:** All results for benzoic acid were estimated (UJ) based on a low laboratory control sample (LCS) recovery.

**Pesticides:** The reporting limits for delta-BHC in Sample Q2-D-3.5 and delta-BHC, aldrin, and gamma chlordane in Sample Q2-D-5.0 were elevated due to matrix interference.

**PCB:** The reporting limit for Aroclor 1248 in Sample Q2-D-5.0 was elevated due to matrix interference.

**NWTPH-Dx:** The diesel range organic result for sample Q2-A-1.0 was estimated (J) based on a low MS recovery and poor laboratory duplicate precision.

**NWTPH-G:** There were no quality control issues that resulted in qualification of data.

**BTEX:** There were no quality control issues that resulted in qualification of data.

**Metals:** The copper results for all soil samples were estimated (J) based on a low MS recovery. All zinc results for the soil samples were estimated (J) based on low MS recovery and poor duplicate precision.

**Conventionals:** There were no quality control issues that resulted in qualification of data.

#### Quendall Terminal Transect Q3 Quendall Terminal

This transect consisted of two soil samples.

**SVOC:** There were no quality control issues that resulted in qualification of data.

**NWTPH-Dx:** There were no quality control issues that resulted in qualification of data.

*Metals:* There were no quality control issues that resulted in qualification of data.

# **Quendall Terminal Borings**

A total of 59 soils and five groundwaters were submitted for the boring sampling. Three sets of field replicates were submitted for the soils: Q4-15.0 & SD-3, Q13-22 & SD-4, and Q15-16.0 & SD-5. One set of field replicates, Q12-W and WD-2, was submitted for the groundwaters.

**PAH:** The water samples required re-extraction as there were no recoveries for the laboratory control sample/laboratory control sample duplicate (LCS/LCSD). Samples Q4-W, Q12-W, Q14-W, Q17-W, and WD-2 were all re-extracted past the holding time. All results for these samples were estimated (J/UJ). There was insufficient sample remaining to re-extract sample Q9-W. All results for this sample were estimated (J) based on the LCS/LCSD recoveries.

For field replicates Q4-15.0 & SD-3, the dibenzofuran results were estimated (J) based on a precision outlier. For field replicates Q12-W and WD-2, the results for acenaphthene, naphthalene, and pyrene were estimated (J).

**NWTPH-Dx:** The diesel range organics and motor oil results were estimated (UJ) based on a surrogate recovery outlier.

**NWTPH-G:** There were no quality control issues that resulted in qualification of data.

**BTEX:** The benzene results for field replicates Q4-15.0 & SD-3 were estimated (J) based on a precision outlier.

**Metals:** The laboratory analyzed the groundwater samples by a method that resulted in much higher reporting limits than called for in the QAPP. The samples were re-analyzed by an alternate method in order to achieve the necessary reporting limits.

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# SAMPLE INDEX POS BNSF ROW Baxter Mills - Transect B1

Sample ID	Lab Sample ID	Matrix	SVOC	TPH-DX	Metals
B1-A-1.0	08-21932-NM59A	SOIL	Χ	Х	Χ
B1-A-3.5	08-21933-NM59B	SOIL	Χ	X	X
B1-A-5.0	08-21934-NM59C	SOIL	X	X	X
B1-A-10.0	08-21935-NM59D	SOIL	X		
B1-A-15.0	08-21936-NM59E	SOIL	Χ		
B1-A-20.0	08-21937-NM59F	SOIL	Χ	X	X
B1-A-24.0	08-21938-NM59G	SOIL	X	Х	Χ
B1-B-1.0	08-21939-NM59H	SOIL	Χ	X	X
B1-B-3.0	08-21940-NM59I	SOIL	X	X	Χ
B1-B-5.0	08-21941-NM59J	SOIL	Χ	Х	Χ
B1-C-5.0	08-21942-NM59K	SOIL	X	X	X
DS-1	08-21943-NM59L	SOIL	Х	X	Х

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Baxter Mill Transect B1 Semivolatile Organic Compounds by Method SW8270D

This report documents the review of analytical data from the analyses of soil samples and the associated laboratory and field quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All samples received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM59	12 Soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

The Chain of Custody (COC) indicated that Samples B1-A-10.0, B1-A-15.0 and B1-B-5.0 were to be placed on hold. However, these samples were analyzed and data reported. No action was taken, other than to note the discrepancy.

#### II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%).

# III. TECHNICAL DATA VALIDATION

The OC requirements that were reviewed are listed below.

1 Holding Times and Sample Preservation

GC/MS Instrument Performance Check Initial Calibration (ICAL)

2 Continuing Calibration (CCAL)

Laboratory Blanks

Field Blanks

Surrogate Compounds

Laboratory Control Samples (LCS/LCSD)

Matrix Spikes/Matrix Spike Duplicates (MS/MSD)

1 Field Replicates

Internal Standards

Target Analyte List

Reporting Limits

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2 Compound Identification and Reported Results

1 Calculation Verification

Ouality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $1.8^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

# **Continuing Calibration**

All relative response factor (RRF) values were acceptable. The continuing calibration (CCAL) percent difference (%D) values were within the ±25% control limits, with the exceptions noted below. When the %D indicated a low bias, positive results and reporting limits were estimated (J/UJ-5B). When the %D outlier indicated a high bias, positive results were estimated (J-5B). No action was taken for non-detects. The following outliers were noted:

**SDG NM59:** CCAL 9/06/08: benzyl alcohol (low bias); 2,4-dinitrophenol (high bias)

CCAL 9/11/08: 2,4-dinitrophenol (high bias)

# Field Replicates

The following acceptance criteria were used to evaluate field precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL (soils) or the RL (waters).

One set of field replicates, B1-A-1.0 and DS-1, was submitted with this SDG. All field precision criteria were met.

# Compound Identification and Reported Results

Samples were screened in order to determine necessary dilutions. In some cases, one or more analytes still exceeded the instrument linear range and required additional dilution. The results for the compounds in the original analyses that were greater than the calibration range were labeled do-not-report (DNR-20). Results for all other compounds in additional dilutions were labeled do not report (DNR-11). The samples requiring additional dilutions are listed below:

**SDG NM59:** B1-B-1.0 (20x)

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were found.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, laboratory control sample, and matrix spike/matrix spike duplicate (MS/MSD) recoveries. Precision was also acceptable as demonstrated by the field replicate and MS/MSD RPD values.

Data were qualified as estimated based on CCAL %D outliers.

Data were labeled do-not-report (DNR) to indicate which result, from multiple dilutions and analyses, should be used. A usable result remains for all analytes in all samples.

Data labeled DNR should not be used for any purpose. All other data, as qualified, are acceptable for use.

B1 SVOC - 3

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Baxter Mill Transect B1 Diesel and Residual Range Hydrocarbons by NWTPH-Dx

This report documents the review of analytical data from the analyses soil samples and the associated field and laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM59	9 Soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

NWTPH-Dx results were reported for the following samples, although no analyses were requested on the chain of custody: B1-A-5.0, B1-B-1.0, B1-B-3.0, and B1-C-5.0.

# II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

# III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are summarized in the table below.

1 Holding Times and Sample Receipt Initial Calibration (ICAL) Continuing Calibration (CCAL) Blanks (Method and Field) Surrogate Compounds Laboratory Control Samples (LCS/LCSD) Matrix Spike/Matrix Spike Duplicates (MS/MSD)

Field Replicates
 Reporting Limits
 Compound Identification
 Sample Results

 Calculation Verification

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Holding Times and Sample Receipt**

Several coolers were received at temperatures that were outside of the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $11.6^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

# Field Replicates

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

**SDG NM59:** One set of field replicates, samples B1-A-1.0 and DS-1, were submitted with this SDG. The RPD for diesel (56.0%) exceeded the control limit. The diesel results for these two samples were estimated (J-9).

# Calculation Verification

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, laboratory control sample/laboratory control sample duplicate (LCS/LCSD), and matrix spike/matrix spike duplicate (MS/MSD) recoveries. Precision was also acceptable as demonstrated by the LCS/LCSD, MS/MSD, and field replicate RPD values, except as noted above.

Data were estimated based on a field replicate RPD outlier.

All data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT

# Pinnacle GeoSciences

# POS BNSF ROW - Baxter Mill Transect B1 Metals by Methods SW6010B, EPA 200.8, and SW7471A and Total Solids by Method 160.3

This report documents the review of analytical data from the analyses of soil samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	 Number of Samples
NM59	9 Soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables, with the exception noted below. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

# II. EDD TO HARDCOPY VERIFICATION

A verification of the electronic data deliverables (EDD) results was performed by comparison to the hardcopy laboratory data package. No errors were found.

# III. TECHNICAL DATA VALIDATION

The QC requirements for review are listed below.

1 Holding Times and Sample Preservation

Initial Calibration

Calibration Verification

**CRDL Standards** 

Laboratory Blanks

Laboratory Control Samples (LCS)

Matrix Spikes (MS)

Laboratory Duplicates

Field Blanks

Field Replicates

Interference Check Samples

Serial Dilutions

ICPMS Internal Standards

Reported Results

1 Calculation Verification

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $1.8^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# Field Replicates

The following acceptance criteria were used to evaluate field precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL (soils) or the RL (waters).

One set of field replicates, Samples B1-A-1.0 and DS-1, was submitted with this SDG. All field precision criteria were met.

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As determined by this evaluation, the laboratory followed the specified analytical methods. The laboratory and field replicate RPD values indicated acceptable precision. Accuracy was also acceptable, as demonstrated by the matrix spike and laboratory control sample recoveries.

No data were qualified for any reason. All data, as reported, are acceptable for use.

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# SAMPLE INDEX POS BNSF ROW

# Quendall Terminal North -Transect Q1

							TPH-G			
Sample ID	Lab Sample ID	Matrix	SVOC	Pest	PCB	TPH-Dx	BTEX	Herbicides	Metals	Turbidity
Q1-A-1.0	08-21754-NM43A	SOIL	X			X			Х	
Q1-A-5.0	08-21755-NM43B	SOIL	Х						X	
Q1-B-1.0	08-21756-NM43C	SOIL	X			Х			Х	
Q1-B-5.0	08-21757-NM43D	SOIL	X						X	
Q1-C-2.5	08-21758-NM43E	SOIL	X			Х			Х	
Q1-C-5.0	08-21759-NM43F	SOIL	X			X			X	, i
Q1-D-2.0	08-21760-NM43G	SOIL	""	Χ	Х				X	
Q1-D-3.5	08-21761-NM43H	SOIL	Х	Χ	Х				X	
Q1-D-5.0	08-21762-NM43I	SOIL	Х						X	
Q1-D-9.0	08-21763-NM43J	SOIL	Х			Х	Х		X	
Q1-D-15.0	08-21764-NM43K	SOIL	X			X	Х		Х	
Q1-D-23.0	08-21765-NM43L	SOIL	Х			Х	Х		X	
Q1-D-30.0	08-21766-NM43M	SOIL	Х			Х	Χ		Х	
SD-2	08-21767-NM43N	SOIL	Х			Х	Χ		X	
Q1-D-W	08-21768-NM43O	WATER	Х			Х	X		X	Х
Q1-Sed	08-21769-NM43P	SED	Х			Х		X	X	
TRIP BLANK	08-21770-NM43Q	WATER					Χ			

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal North Transect Q1 Semivolatile Organic Compounds by Method SW8270D

This report documents the review of analytical data from the analyses of groundwater, soil, and sediment samples and the associated laboratory and field quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All samples received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM43	1 Groundwater, 13 soil, 1 sediment

#### I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

# II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%).

# III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are listed below.

- Holding Times and Sample Preservation
   GC/MS Instrument Performance Check
- 1 Initial Calibration (ICAL) Continuing Calibration (CCAL) Laboratory Blanks Field Blanks
  - Surrogate Compounds
- 1 Laboratory Control Samples (LCS/LCSD)

- 2 Matrix Spikes/Matrix Spike Duplicates (MS/MSD)
- Field Replicates
   Internal Standards
   Target Analyte List
   Reporting Limits
- 2 Compound Identification and Reported Results
- 1 Calculation Verification

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $1.8^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

#### **Initial Calibration**

All relative response factor (RRF) values were greater than the 0.05 minimum control limit. The initial calibration (ICAL) percent relative standard deviation (%RSD) values were within the  $\pm 30\%$  control limits, with the exceptions noted below. Positive results associated with %RSD outliers were estimated (J-5A). No action was taken for non-detects. The following outliers were noted:

**SDG NM43:** ICAL 9/12/08: 2,4-dinitrophenol (31.2%). This analyte was not detected in the associated samples, therefore no qualification of data was necessary.

# **Laboratory Control Samples**

Laboratory control sample/laboratory control sample duplicates (LCS/LCSD) analyses were performed at the required frequency. The recovery values were within the specified control limits, with the exceptions noted below.

The LCSD %R value for hexachloroethane was less than the lower control limit. The LCS %R was acceptable, therefore no qualification of data was necessary. The relative percent difference (RPD) values for hexachloroethane and hexachlorobutadiene were greater than the control limit. No positive results were associated with these outliers; no qualification was necessary.

# Matrix Spike/Matrix Spike Duplicates

Matrix spike/matrix spike duplicate (MS/MSD) analyses were performed at the required frequency. The MS/MSD recovery values were within the specified control limits, with the exceptions noted below. If the outliers indicated a potential high bias, associated positive results in the parent sample only were qualified as estimated (J-8). If the outliers indicated a potential low bias, positive results and reporting limits in the parent sample were estimated (J/UJ-8).

The MS/MSD RPD values were within the specified control limits, with the exceptions noted below. For RPD outliers, positive results only in parent sample are estimated (J-9). No action is taken for non-detects.

The following outliers were noted:

Sample Q1-D-3.5 MS/MSD: Benzoic acid was not detected in the MS. The MSD %R was acceptable. The MS/MSD RPD value was greater than the control limit. The benzoic acid result for the parent sample was estimated (UJ-8).

# Field Replicates

The following acceptance criteria were used to evaluate field precision: the RPD control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL (soils) or the RL (waters).

**SDG NM43:** One set of field replicates, Q1-D-23.0 and SD-2, was submitted with this SDG. The following RPD outliers were noted:

benzo(a)anthracene	57.4%	benzo(a)pyrene	53.3%
benzo(b)fluoranthene	58.1%	benzo(g,h,i)perylene	55.2%
chrysene	53.8%	fluoranthene	51.9%
indeno(1,2,3-cd)pyrene	71.2%	pyrene	50.6%

The results for these analytes in the above samples were estimated (J-9).

# Compound Identification and Reported Results

Samples were screened in order to determine necessary dilutions. In some cases, one or more analytes still exceeded the instrument linear range and required additional dilution. The results for the compounds in the original analyses that were greater than the calibration range were labeled do-not-report (DNR-20). Results for all other compounds in additional dilutions were labeled do not report (DNR-11). The samples requiring additional dilutions are:

Q1-A-1.0 (10x)	Q1-B-1.0 (25x)	Q1-D-9.0 (30x)
Q1-D-15.0 (30x)	Q1-D-23.0 (30x)	SD-2 (30x)

#### **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were found.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, LCS/LCSD, and MS/MSD %R values, with the exceptions noted above. Precision was also acceptable as demonstrated by the field replicate, MS/MSD, and LCS/LCSD RPD values.

Data were qualified as estimated based on CCAL %D and field replicate RPD outliers.

Data were labeled do-not-report (DNR) to indicate which result, from multiple dilutions and analyses, should be used. A usable result remains for all analytes in all samples.

Data labeled DNR should not be used for any purpose. All other data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal North Transect Q1 Chlorinated Pesticides - EPA Method 8081A

This report documents the review of analytical data from the analyses of soil samples and the associated laboratory quality control (QC) samples. Samples were analyzed by Analytical Resources, Inc., Tukwila, Washington. All data received a full (Level IV) validation. Please see the Sample Index for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM43	2 Soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

#### II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

#### III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are listed below.

1 Holding Times and Sample Preservation

Initial Calibration (ICAL)

Continuing Calibration (CCAL)

Laboratory Blanks

Field Blanks

Surrogate Compounds

Laboratory Control Samples (LCS/LCSD)

Matrix Spikes/Matrix Spike Duplicates (MS/MSD)

Field Replicates

Internal Standards

Compound Identification

Reporting Limits

I Calculation Verification

# **Holding Times and Sample Preservation**

The temperatures of several coolers were outside the recommended temperature range of  $4^{\circ}\text{C} \pm 2^{\circ}$  when received at the laboratory, ranging from 6.8° to 11.6°C. The temperature outliers did not impact data quality and no data were qualified.

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory performed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, matrix spike/matrix spike duplicate (MS/MSD), and laboratory control sample/laboratory control sample duplicate (LCS/LCSD) recoveries. Precision was acceptable as demonstrated by the relative percent difference values for the LCS/LCSD and MS/MSD analyses.

No data were qualified for any reason. All data, as reported, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal North Transect Q1 PCB Aroclors by SW846 Method 8082

This report documents the review of analytical data from the analyses of soil samples and the associated laboratory and field quality control (QC) samples. Samples were analyzed by Analytical Resources, Inc., Tukwila, Washington. All data received a full (Level IV) validation. Please see the Sample Index for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM43	2 Soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

# II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

# III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are listed below.

1 Holding Times and Sample Preservation

Initial Calibration (ICAL)

Continuing Calibration (CCAL)

Laboratory Blanks

Field Blanks

Surrogate Compounds

Laboratory Control Samples (LCS/LCSD)

Matrix Spikes/Matrix Spike Duplicates (MS/MSD)

Field Replicates

Internal Standards

Compound Identification

Reporting Limits

Calculation Verification

# **Holding Times and Sample Preservation**

The temperatures of several coolers were outside the recommended temperature range of  $4^{\circ}\text{C} \pm 2^{\circ}$  when received at the laboratory, ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . The temperature outliers did not impact data quality and no data were qualified.

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory performed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, matrix spike/matrix spike duplicate (MS/MSD), and laboratory control sample/laboratory control sample duplicate (LCS/LCSD) percent recovery values. Precision was also acceptable as demonstrated by the relative percent difference values for the LCS/LCSD and MS/MSD analyses.

No data were qualified for any reason. All data, as reported, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal North Transect Q1 Herbicides by EPA Method 8151A

This report documents the review of analytical data from the analyses of one soil sample and the associated laboratory quality control (QC) samples. The sample was analyzed by Analytical Resources, Inc., Tukwila, Washington. All data received a full (Level IV) validation. Please see the Sample Index for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM43	1 Soil

#### I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

#### II. **EDD TO HARDCOPY VERIFICATION**

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%).

#### III. **TECHNICAL DATA VALIDATION**

The QC requirements that were reviewed are listed below.

- 1 Holding Times and Sample Preservation Instrument Breakdown Check Initial Calibration (ICAL) Continuing Calibration (CCAL) Laboratory Blanks
- 1 Surrogate Compounds
- 2 Laboratory Control Samples (LCS/LCSD)
- 2 Matrix Spikes/Matrix Spike Duplicates (MS/MSD) Field Replicates Target Analyte List
  - Reporting Limits
  - Compound Identification
- Calculation Verification

# **Holding Times and Sample Preservation**

The temperatures of several coolers were outside the recommended temperature range of 4°C ±2° when received at the laboratory, ranging from 6.8° to 11.6°C. The temperature outliers did not impact data quality and no data were qualified.

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Ouglity control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below,

# **Surrogate Compounds**

The recovery for the surrogate compound 2,4-dichlorophenylacetic acid (DCPA) was less than the lower control limit in QC sample LCS-090208. No qualifiers are assigned to QC samples.

# **Laboratory Control Samples**

Laboratory control sample/laboratory control sample duplicates (LCS/LCSD) were analyzed. The LCS recoveries for five analytes were less than the lower control limits. The LCSD recoveries were acceptable, therefore no action was taken. The relative percent difference (RPD) values for these five analytes were greater than the control limits. These analytes were not detected in the associated sample, therefore no qualifiers were required.

The LCS/LCSD recoveries for dinoseb were less than 10%. This analyte was not detected in the associated sample; the reporting limit was rejected (R-10)

# Matrix Spike/Matrix Spike Duplicates

Matrix spike/matrix spike duplicate (MS/MSD) analyses were performed at the required frequency. The MS/MSD percent recovery (%R) values were within the specified control limits, with the following exceptions:

The %R value for dinoseb in the MS was less than the lower control limit. Dinoseb was not recovered in the MSD. The result for this compound was rejected (R-8) in the parent sample.

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory performed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, MS/MSD, and LCS/LCSD %R values, except as noted above. Precision was acceptable as demonstrated by the RPD values for the LCS/LCSD and MS/MSD analyses.

The dinoseb result was rejected based on MS/MSD and LCS/LCSD recoveries that were less than 10%.

Rejected data should not be used for any purpose. All other data, as reported, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal North Transect Q1 Diesel and Residual Range Hydrocarbons by NWTPH-Dx

This report documents the review of analytical data from the analyses of groundwater, soil, and sediment samples and the associated field and laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM43	1 Groundwater, 9 soil, 1 sediment

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

# II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

# III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are summarized in the table below.

 Holding Times and Sample Preservation Initial Calibration (ICAL)
 Continuing Calibration (CCAL)
 Blanks (Method and Field)

Surrogate Compounds
 Laboratory Control Samples (LCS/LCSD)

Matrix Spike/Matrix Spike Duplicates (MS/MSD)

Field Replicates
 Reporting Limits
 Compound Identification
 Sample Results

 Calculation Verification

Quality control results are discussed below, but no data were qualified.

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures that were outside of the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $11.6^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Surrogate Compounds**

The surrogate recoveries for several samples could not be calculated due to necessary dilutions. No action was necessary.

# **Field Replicates**

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

One set of field replicates, samples Q1-D-23.0 and SD-2 were submitted with this SDG. The RPD values for diesel (53.7%) and motor oil (51.0%) exceeded the control limit. The results for these analytes were estimated (J-9) in both samples.

#### **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, laboratory control sample/laboratory control sample duplicate (LCS/LCSD), and matrix spike/matrix spike duplicate (MS/MSD) recoveries. Precision was also acceptable as demonstrated by the LCS/LCSD, MS/MSD, and field replicate RPD values, except as noted above.

Data were estimated because of a field duplicate RPD outlier.

All data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal North Transect Q1 Gasoline Range Organics by NWTPH-Gx

This report documents the review of analytical data from the analyses of groundwater and soil samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM43	1 Groundwater, 5 soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

# II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found, with the following exception:

# III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are summarized in the table below.

1	Holding Times and Sample Receipt		Laboratory Duplicates
	Instrument Performance Check		Laboratory Control Samples
	Initial Calibration (ICAL)	1	Field Replicates
	Continuing Calibration (CCAL)		Reported Results
	Blanks (Method and Field)		Compound Identification
	Surrogate Compounds	1	Calculation Verification (full validation of

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

# **Holding Times and Sample Receipt**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $1.8^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

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<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Field Replicates**

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

One set of field replicates, Samples Q1-D-23.0 and SD-2 was submitted. All field precision criteria were met.

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate and laboratory control sample percent recovery values. Precision was acceptable as demonstrated by the RPD values for the laboratory and field replicate analyses.

No data were qualified for any reason. All data, as reported, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal North Transect Q1 BTEX Compounds by 8021Mod

This report documents the review of analytical data from the analyses of groundwater and soil samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM43	1 Groundwater, 5 soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

# II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

#### III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are summarized in the table below.

 Holding Times and Sample Preservation Instrument Performance Check Initial Calibration (ICAL)
 Continuing Calibration (CCAL)
 Blanks (Method and Field)

Surrogate Compounds

Laboratory Control Samples (LCS/LCSD)

Field Replicates
 Reported Results
 Compound Identification

 Calculation Verification

<sup>1</sup> Quality control results are discussed below, but no data were qualified.

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures outside of the recommended range of  $4^{\circ}$ C  $\pm 2^{\circ}$ , ranging from 1.6° to 11.6°C. It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Field Replicates**

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

One set of field replicates, Samples Q1-D-23.0 and SD-2 was submitted with this SDG. All field precision criteria were met.

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate and laboratory control sample/laboratory control sample duplicate (LCS/LCSD) percent recovery values. Precision was acceptable as demonstrated by the LCS/LCSD and field replicate RPD values.

No data were qualified for any reason. All data, as reported, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal North Transect Q1 Metals by Methods SW6010B, EPA 200.8, and SW7471A

This report documents the review of analytical data from the analyses of groundwater, soil, and sediment samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples	
NM43	1 Groundwater, 14 soil, 1 sediment	

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables, with the exception noted below. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

The nickel results were not reported for the QC samples, although this was a target analyte for some of the samples in the preparation batch. Results were calculated from the raw data and no further action was taken.

# II. EDD TO HARDCOPY VERIFICATION

A verification of the electronic data deliverables (EDD) results was performed by comparison to the hardcopy laboratory data package. No errors were found.

#### III. TECHNICAL DATA VALIDATION

The QC requirements for review are listed below.

 Holding Times and Sample Preservation Initial Calibration
 Calibration Verification
 CRDL Standards
 Laboratory Blanks

Laboratory Control Samples (LCS)

Matrix Spikes (MS)

Laboratory Duplicates

Field Replicates
Interference Check Samples

interierence Offeck Samples

Serial Dilutions

ICPMS Internal Standards

Reported Results

Calculation Verification

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

recommended range of 4°C ±2°, ranging from 1.6° to 1.8°C. It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

# Laboratory Blanks

Zinc was detected in the method blank at levels greater than the method detection limit (MDL). To evaluate the effect on the sample data, an action level of five times the blank concentration was established. All associated sample results were greater than the action level, therefore no qualification of the data was necessary.

# **Field Replicates**

The following acceptance criteria were used to evaluate field precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL (soils) or the RL (waters).

One set of field replicates, Samples Q1-D-23.0 and SD-2 was submitted. There were no positive results in either sample. Field precision was acceptable.

#### **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As determined by this evaluation, the laboratory followed the specified analytical methods. The laboratory and field replicate RPD values indicated acceptable precision. Accuracy was also acceptable, as demonstrated by the matrix spike and laboratory control sample recoveries.

No data were qualified for any reason. All data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT

# **Pinnacle GeoSciences**

# POS BNSF ROW - Quendall Terminal North Transect Q1 Turbidity by Method 180.1 and Total Solids by Method 160.3

This report documents the review of analytical data from the analyses of groundwater, sediment, and soil samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples	
NM43	1 Groundwater, 14 soil, 1 sediment	

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

# II. EDD TO HARDCOPY VERIFICATION

A verification of the electronic data deliverables (EDD) results was performed by comparison to the hardcopy laboratory data package. No errors were found.

#### III. TECHNICAL DATA VALIDATION

The QC requirements for review are listed below.

1 Holding Times and Sample Preservation

Initial Calibration

Calibration Verification

1 Field Replicates

Laboratory Blanks

Laboratory Control Samples (LCS)

Matrix Spike (MS)

Laboratory Replicates

Reporting Limits

Calculation Verification

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $1.8^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# Field Replicates

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

No data were qualified based on field replicate precision outliers. Users of the data should consider the impact of field precision outliers on the reported results.

One set of field replicates, Samples Q1-D-23.0 and SD-2 was submitted with this SDG. All field precision criteria were met.

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As determined by this evaluation, the laboratory followed the specified analytical methods. The laboratory and field replicate RPD values indicated acceptable precision. Accuracy was also acceptable, as demonstrated by the laboratory control sample recoveries.

No data were qualified for any reason. All data, as reported, are acceptable for use.

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# SAMPLE INDEX POS BNSF ROW

# **Quendall Terminal South - Transect Q2**

		1					TPH-G		
Sample ID	Lab Sample ID	Matrix	svoc	Pest	PCB	TPH-Dx	BTEX	Metals	Turbidity
Q2-A-1.0	08-21714-NM40A	SOIL	Χ			Х		Χ	
Q2-A-5.0	08-21715-NM40B	SOIL	Χ			X		Х	
Q2-B-1.0	08-21716-NM40C	SOIL	Х			Х		Х	
Q2-B-5.0	08-21717-NM40D	SOIL	Х					X	
Q2-C-3.5	08-21718-NM40E	SOIL	Х			Х		Χ	
Q2-C-5.0	08-21719-NM40F	SOIL	Х			Х		X	
Q2-C-13.0	08-21720-NM40G	SOIL	Х			Х	Х	Χ	
Q2-C-25.0	08-21721-NM40H	SOIL	Х						
Q2-D-3.5	08-21722-NM40I	SOIL	Х	X	Х	Х		Х	
Q2-D-5.0	08-21723-NM40J	SOIL	Х	Х	Х	Х		Х	
Q2-D-10.0	08-21724-NM40K	SOIL	Χ			X		X	
Q2-D-13.0	08-21725-NM40L	SOIL	Χ			Х	Х	Χ	
Q2-D-18.0	08-21726-NM40M	SOIL	Х			X	Х	X	
Q2-D-35.0	08-21727-NM40N	SOIL	Χ			X	Х	Х	
Q2-D-W	08-21728-NM40O	WATER					Х	X	Χ
WD-1	08-21729-NM40P	WATER					Х	Х	Χ
TRIP BLANK	08-21730-NM40Q	WATER					X		

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal South Transect Q2 Semivolatile Organic Compounds by Method SW8270D

This report documents the review of analytical data from the analyses of soil samples and the associated laboratory and field quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All samples received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM40	14 Soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

Water samples Q2-D-W and WD-1 appeared on the Chain of Custody (COC), but were not analyzed by the laboratory.

#### II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%).

#### III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are listed below.

- Holding Times and Sample Preservation GC/MS Instrument Performance Check Initial Calibration (ICAL)
- Continuing Calibration (CCAL)
   Laboratory Blanks
   Field Blanks
  - Surrogate Compounds
- 2 Laboratory Control Samples (LCS/LCSD)

- 1 Matrix Spikes/Matrix Spike Duplicates (MS/MSD)
  - Field Replicates
  - Internal Standards
  - Target Analyte List
  - Reporting Limits
- 2 Compound Identification and Reported Results
- 1 Calculation Verification

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

recommended range of 4°C  $\pm$ 2°, ranging from 1.6° to 1.8°C. It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

# **Continuing Calibration**

All relative response factor (RRF) values were greater than the 0.05 minimum control limit. The continuing calibration (CCAL) percent difference (%D) values were within the ±25% control limits, with the exceptions noted below. When the %D indicated a low bias, positive results and reporting limits were estimated (J/UJ-5B). When the %D outlier indicated a high bias, positive results were estimated (J-5B). No action was taken for non-detects. The following outliers were noted:

CCAL 9/17/08: hexachlorocyclopentadiene, 2,4-dinitrophenol, bis(2-ethylhexyl) phthalate (high bias)

# **Surrogates**

The percent recovery (%R) values for the surrogates were within the specified control limits with the exception noted below. Several samples required dilutions greater than 10x. Surrogate recovery control limits do not apply in these cases.

The %R value for d14-p-terphenyl was less than the lower control limit in Sample Q2-D-5.0. All other surrogate recoveries were in control. One outlier per acid or base/neutral fraction is allowed, therefore no qualification was necessary for the single outlier.

# **Laboratory Control Samples**

Laboratory control sample/laboratory control sample duplicates (LCS/LCSD) analyses were performed at the required frequency. The recovery values were within the specified control limits, with the exceptions noted below.

The LCS/LCSD R% values for benzoic acid were less than the lower control limit. Benzoic acid was not detected in the associated samples; reporting limits were estimated (UJ-10).

# Matrix Spike/Matrix Spike Duplicates

Matrix spike/matrix spike duplicate (MS/MSD) analyses were performed at the required frequency. The MS/MSD recovery values were within the specified control limits, with the exceptions noted below. If the outliers indicated a potential high bias, associated positive results in the parent sample only were qualified as estimated (J-8). If the outliers indicated a potential low bias, positive results and reporting limits in the parent sample were estimated (J/UJ-8).

The MS/MSD relative percent difference (RPD) values were within the specified control limits, with the exceptions noted below. For RPD outliers, positive results only in parent sample are estimated (J-9). No action is taken for non-detects.

One outlier was noted, QC Sample Q2-A-5.0, the MSD %R value for chrysene was less than the lower control limit. No qualification was necessary as the MS %R was in control.

# **Compound Identification and Reported Results**

Samples were screened in order to determine necessary dilutions. In some cases, one or more analytes still exceeded the instrument linear range and required additional dilution. The results for the compounds in the original analyses that were greater than the calibration range were labeled do-not-report (DNR-20). Results for all other compounds in additional dilutions were labeled do not report (DNR-11). The samples requiring additional dilutions are listed below:

**SDG NM40:** Q2-C-13.0 (20x), Q2-C-3.5 (5x), Q2-D-5.0 (100x), Q2-D-10.0 (40x), Q2-D-13.0 (5x), Q2-D-18.0 (40x), Q2-D-3.5 (30x)

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were found.

### IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, LCS/LCSD, and MS/MSD %R values, with the exceptions noted above. Precision was also acceptable as demonstrated by the field replicate, MS/MSD, and LCS/LCSD RPD values.

Data were qualified as estimated based on LCS/LCSD %R outliers.

Data were labeled do-not-report (DNR) to indicate which result, from multiple dilutions and analyses, should be used. A usable result remains for all analytes in all samples.

Data labeled DNR should not be used for any purpose. All other data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal South Transect Q2 Chlorinated Pesticides - EPA Method 8081A

This report documents the review of analytical data from the analyses of soil samples and the associated laboratory quality control (QC) samples. Samples were analyzed by Analytical Resources, Inc., Tukwila, Washington. All data received a full (Level IV) validation. Please see the Sample Index for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM40	2 Soil

### I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

### II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

#### III. TECHNICAL DATA VALIDATION

The OC requirements that were reviewed are listed below.

- 1 Holding Times and Sample Preservation Initial Calibration (ICAL)
  - Continuing Calibration (CCAL)

Laboratory Blanks

Field Blanks

1 Surrogate Compounds

Laboratory Control Samples (LCS/LCSD)

- 1 Matrix Spikes/Matrix Spike Duplicates (MS/MSD) Field Replicates
- 1 Internal Standards
- 2 Compound Identification
- 2 Reporting Limits
- 1 Calculation Verification

# **Holding Times and Sample Preservation**

The temperatures of several coolers were outside the recommended temperature range of  $4^{\circ}\text{C} \pm 2^{\circ}$  when received at the laboratory, ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . The temperature outliers did not impact data quality and no data were qualified.

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

Surrogate Compounds

The surrogate percent recovery (%R) values were acceptable, with the exceptions noted below.

The samples were analyzed at dilutions (10x and 100x), therefore no surrogate recoveries were reported. No action was taken on this basis.

Matrix Spike/Matrix Spike Duplicates

Matrix spike/matrix spike duplicate (MS/MSD) analyses were not performed. Accuracy and precision were assessed using the laboratory control sample and laboratory control sample duplicate (LCS/LCSD).

Internal Standards

The recoveries of the internal standard hexabromobiphenyl were greater than the 200% control limit for Samples Q2-D-3.5 and Q2-D-5.0. There were no positive results in these samples, therefore no qualifiers were assigned.

Reporting Limits

Reporting limits were elevated when chromatograms indicated non-target background interference. The reporting limits for these analytes were flagged "Y" by the laboratory. These "Y" flagged results were qualified as not-detected (U-22). The following samples had elevated reporting limits for one or more compounds:

**SDG NM40:** Sample Q2-D-3.5: delta-BHC

Sample Q2-D-5.0: delta-BHC, aldrin, and gamma chlordane

**Compound Identification and Reported Results** 

Samples Q2-D-3.5 and Q2-D-5.0 were re-analyzed at a 100x dilution because of internal standard outliers in the initial 10x dilutions. No positive results were reported for either analysis. The dilution analyses were labeled as do-not-report (DNR-11) in favor of the initial analyses.

Calculation Verification

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory performed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, MS/MSD, and LCS/LCSD %R values. Precision was acceptable as demonstrated by the relative percent difference values for the LCS/LCSD and MS/MSD analyses.

Data were qualified as not detected at elevated reporting limits due to matrix interferences.

Data were labeled as do-not-report (DNR) to indicate which set of data from multiple analyses should be used. A usable result for remains for every analyte in each sample, therefore completeness is unaffected.

Data labeled DNR should not be used for any purpose. All other data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal South Transect Q2 PCB Aroclors by SW846 Method 8082

This report documents the review of analytical data from the analyses of soil samples and the associated laboratory and field quality control (QC) samples. Samples were analyzed by Analytical Resources, Inc., Tukwila, Washington. All data received a full (Level IV) validation. Please see the Sample Index for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM40	2 Soil

## I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

# II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

### III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are listed below.

1 Holding Times and Sample Preservation

Initial Calibration (ICAL)

Continuing Calibration (CCAL)

Laboratory Blanks

Field Blanks

Surrogate Compounds

Laboratory Control Samples (LCS/LCSD)

1 Matrix Spikes/Matrix Spike Duplicates (MS/MSD)

Field Replicates

Internal Standards

Compound Identification

2 Reporting Limits

1 Calculation Verification

# **Holding Times and Sample Preservation**

The temperatures of several coolers were outside the recommended temperature range of  $4^{\circ}\text{C} \pm 2^{\circ}$  when received at the laboratory, ranging from 6.8° to 11.6°C. The temperature outliers did not impact data quality and no data were qualified.

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# Matrix Spike/Matrix Spike Duplicates

Matrix spike/matrix spike duplicate (MS/MSD) analyses were performed. Accuracy and precision were assessed using the laboratory control sample and laboratory control sample duplicate (LCS/LCSD).

# **Reporting Limits**

Reporting limits were elevated when chromatograms indicated non-target background interference. The reporting limits for these analytes were flagged "Y" by the laboratory. These "Y" flagged results were qualified as not-detected (U-22). The following samples had elevated reporting limits for one or more compounds:

**SDG NM40:** Sample Q2-D-5.0: Aroclor 1248

#### **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

### IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory performed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, MS/MSD, and LCS/LCSD percent recovery values. Precision was also acceptable as demonstrated by the relative percent difference values for the LCS/LCSD and MS/MSD analyses.

Reporting limits were elevated based on matrix interference.

All data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal South Transect Q2 Diesel and Residual Range Hydrocarbons by NWTPH-Dx

This report documents the review of analytical data from the analyses of soil samples and the associated field and laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM40	12 Soil

#### I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

Water samples Q2-D-W and WD-1 appeared on the Chain of Custody (COC), but were not analyzed by the laboratory.

### II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

#### III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are summarized in the table below.

 Holding Times and Sample Preservation Initial Calibration (ICAL) Continuing Calibration (CCAL) Blanks (Method and Field)
 Surrogate Compounds Laboratory Control Samples (LCS/LCSD) 2 Matrix Spike/Matrix Spike Duplicates (MS/MSD)
Field Replicates
Reporting Limits
Compound Identification
Sample Results

1 Calculation Verification

<sup>1</sup> Quality control results are discussed below, but no data were qualified.

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures that were outside of the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Surrogate Compounds**

The surrogate recoveries for several samples could not be calculated due to necessary dilutions. No action was necessary.

# Matrix Spike/Matrix Spike Duplicates

Matrix spike/matrix spike duplicate samples (MS/MSD) were analyzed at the proper frequency of one per 20 samples or one per batch; whichever was more frequent. The percent recovery (%R) and relative percent difference (RPD) values were within the laboratory control limits, with the exception noted below.

QC Sample Q2-A-1.0: diesel (MSD 9.6 %; RPD 39.6%) – result for the parent sample was estimated (J-8,9).

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

## IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, laboratory control sample/laboratory control sample duplicate (LCS/LCSD), and MS/MSD %R values, except as noted above. Precision was also acceptable as demonstrated by the MS/MSD and LCS/LCSD RPD values, except as noted above.

Data were estimated because of MS/MSD %R and RPD outliers.

All data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal South Transect Q2 Gasoline Range Organics by NWTPH-Gx

This report documents the review of analytical data from the analyses of groundwater and soil samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM40	2 Groundwater, 4 soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

# II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found, with the following exception:

# III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are summarized in the table below.

•		Laboratory Duplicates Laboratory Control Samples (LCS)
n (ICAL)	1	Field Replicates
bration (CCAL)		Reported Results
and Field)		Compound Identification
pounds	1	Calculation Verification
f	and Sample Receipt formance Check on (ICAL) ibration (CCAL) d and Field) upounds	formance Check on (ICAL) 1 ibration (CCAL) d and Field)

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

# **Holding Times and Sample Receipt**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $1.8^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below,

# **Field Replicates**

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

One set of field replicates, Samples Q2-D-W and WD-1 was submitted. All field precision criteria were met.

#### Calculation Verification

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate and laboratory control sample percent recovery values. Precision was acceptable as demonstrated by the RPD values for the laboratory and field replicate analyses.

No data were qualified for any reason. All data, as reported, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal South Transect Q2 BTEX Compounds by EPA Method 8021Mod

This report documents the review of analytical data from the analyses of groundwater and soil samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM40	2 Groundwater, 4 soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

# II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

# III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are summarized in the table below.

1 Holding Times and Sample Preservation

Instrument Performance Check

Initial Calibration (ICAL)

Continuing Calibration (CCAL)
Blanks (Method and Field)

. . .

Surrogate Compounds

Laboratory Control Samples (LCS/LCSD)

1 Field Replicates

Reported Results

Compound Identification

1 Calculation Verification

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures outside of the recommended range of  $4^{\circ}\text{C}$   $\pm 2^{\circ}$ , ranging from 1.6° to 11.6°C. It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below,

# Field Replicates

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

One set of field replicates, Samples Q2-D-W and WD-1 was submitted. All field precision criteria were met.

### **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

### IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate and laboratory control sample/laboratory control sample duplicate (LCS/LCSD) percent recovery values. Precision was acceptable as demonstrated by the RPD values for the LCS/LCSD and field replicate analyses.

No data were qualified for any reason. All data, as reported, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal South Transect Q2 Metals by Methods SW6010B, EPA 200.8, and SW7471A

This report documents the review of analytical data from the analyses of groundwater and soil samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples	
NM40	2 Groundwater, 13 soil	

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables, with the exception noted below. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

## II. EDD TO HARDCOPY VERIFICATION

A verification of the electronic data deliverables (EDD) results was performed by comparison to the hardcopy laboratory data package. No errors were found.

### III. TECHNICAL DATA VALIDATION

The QC requirements for review are listed below.

- Holding Times and Sample Preservation Initial Calibration
   Calibration Verification
   CRDL Standards
   Laboratory Blanks
   Laboratory Control Samples (LCS)
   Matrix Spikes (MS)
- 2 Laboratory Duplicates
- 1 Field Replicates Interference Check Samples Serial Dilutions ICPMS Internal Standards Reported Results
- 1 Calculation Verification

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of 4°C  $\pm 2^{\circ}$ , ranging from 6.8° to 11.6°C. Several coolers were received at temperatures less than the recommended range of 4°C  $\pm 2^{\circ}$ , ranging from 1.6° to 1.8°C. It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Laboratory Blanks**

Zinc was detected in the method blank at levels greater than the method detection limit (MDL). To evaluate the effect on the sample data, an action level of five times the blank concentration was established. All associated sample results were greater than the action level, therefore no qualification of the data was necessary.

# **Matrix Spikes**

A matrix spike sample (MS) was analyzed at the proper frequency of one per 20 samples or one per batch; whichever was more frequent. The percent recovery (%R) values were within the laboratory control limits, with the exceptions noted below. For %R values less than the lower control limit, the associated positive results and non-detects were estimated (J/UJ-8) to indicate a possible low bias. The following outliers were noted:

QC Sample Q2-D-3.5 (soil): copper (72.6%), zinc (-21.1%) – low bias. Although there was no recovery for zinc in the MS, the results were estimated rather than rejected due to the non-homogenous nature of the sample as indicated by the duplicate results.

# **Laboratory Duplicates**

Laboratory duplicate relative percent difference (RPD) values were used to evaluate precision. The RPD values were within the control limit of 35% for soil sample results greater than five times the reporting limit (for results less than five times the reporting limit, the difference was less than twice the reporting limit) with the exceptions noted below. For RPD or difference values exceeding the control limits, associated positive results and non-detects were estimated (J/UJ-9). The following outliers were noted:

QC Sample Q2-D-3.5 - soil: zinc (37.3%)

# Field Replicates

The following acceptance criteria were used to evaluate field precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL (soils) or the RL (waters).

One set of field replicates, Samples Q2-D-W and WD-1 was submitted with this SDG. All field precision criteria were met.

### **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As determined by this evaluation, the laboratory followed the specified analytical methods. The laboratory and field replicate RPD values indicated acceptable precision, except as noted above. Accuracy was also acceptable, as demonstrated by the MS and laboratory control sample recoveries, except as noted above.

Data were estimated based on MS %R and laboratory duplicate RPD outliers.

All data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal South Transect Q2 Turbidity by Method 180.1 and Total Solids by Method 160.3

This report documents the review of analytical data from the analyses of groundwater and soil samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NM40	2 Groundwater, 13 soil

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

#### II. EDD TO HARDCOPY VERIFICATION

A verification of the electronic data deliverables (EDD) results was performed by comparison to the hardcopy laboratory data package. Errors in the units for the turbidity duplicate analysis for samples Q2-D-W and Q4-W were found. The EDD was corrected and no further action was taken.

#### III. TECHNICAL DATA VALIDATION

The QC requirements for review are listed below.

1	Holding Times and Sample Preservation		Laboratory Control Samples
	Initial Calibration		Matrix Spikes (MS)
	Calibration Verification		Laboratory Replicates
1	Field Replicates		Reporting Limits
	Laboratory Blanks	1	Calculation Verification

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $1.8^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Field Replicates**

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

No data were qualified based on field replicate precision outliers. Users of the data should consider the impact of field precision outliers on the reported results.

One set of field replicates, Samples Q2-D-W and WD-1 was submitted with this SDG. All field precision criteria were met.

#### Calculation Verification

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As determined by this evaluation, the laboratory followed the specified analytical methods. The laboratory replicate RPD values indicated acceptable precision. Accuracy was also acceptable, as demonstrated by the laboratory control sample recoveries.

No data were qualified for any reason. All data, as reported, are acceptable for use.

# SAMPLE INDEX POS BNSF ROW Quendall Terminal - Borings

Sample ID	Lab Sample ID	Matrix	PAH	TPH-Dx	TPH-G BTEX	Metals	Turbidity
Q4-2.5	08-29725-NX66A	SOIL	X	X		X	
Q4-15.0	08-29727-NX66C	SOIL	X	X	Х	X	<u> </u>
Q4-27.0	08-29729-NX66E	SOIL	X	X		Х	
Q4-31.0	08-29730-NX66F	SOIL	X	Х	X	X	
Q5-14.0	08-29732-NX66H	SOIL	X	X	Χ	Х	† <del>***</del>
Q5-18.0	08-29733-NX66I	SOIL	X	X	X	X	
Q5-25.5	08-29734-NX66J	SOIL	X	X		Х	
Q6-4.0	08-29736-NX66K	SOIL	X	X		Х	
Q6-18.0	08-29738-NX66M	SOIL	X	X	Х	X	
Q6-22.5	08-29739-NX66N	SOIL	X	X		Х	
Q7-4.0	08-29740-NX66O	SOIL	X	X		Х	
Q7-5.5	08-29741-NX66P	SOIL	Х	X		X	1
Q7-9.0	08-29742-NX66Q	SOIL	X	X	Х	X	
Q7-19.5	08-29743-NX66R	SOIL	X	X	X	X	
Q8-3.5	08-29744-NX66S	SOIL	X	$\frac{1}{X}$		X	
Q8-16.0	08-29758-NX66T	SOIL	X	$\frac{\hat{x}}{\hat{x}}$	Χ	X	
Q8-24.0	08-29767-NX71A	SOIL	X	$\frac{1}{x}$		X	
Q8-28.0	08-29768-NX71B	SOIL	X	$\frac{1}{x}$		X	
Q9-18.0	08-29770-NX71D	SOIL	$\frac{x}{x}$	X	Х	X	
Q9-25.0	08-29771-NX71E	SOIL	X	X		X	<del> </del>
Q9-28.0	08-29772-NX71F	SOIL	X	$\frac{1}{x}$	Χ	X	
Q10-5.0	08-29773-NX71G	SOIL	$\frac{x}{x}$	$\frac{x}{x}$		X	<del>                                     </del>
Q10-19.0	08-29775-NX71I	SOIL	X	$\frac{1}{x}$		X	
Q10-26.0	08-29776-NX71J	SOIL	$\frac{x}{x}$	$\frac{x}{x}$	Χ	X	
Q11-11.5	08-29778-NX71L	SOIL	X	$\frac{1}{X}$	X	X	<b></b>
Q11-18.0	08-29779-NX71M	SOIL	X	$\frac{1}{X}$		X	
Q11-26.0	08-29780-NX71N	SOIL	X	$\frac{x}{x}$	X	X	<b>!</b>
Q12-4.5	08-29781-NX710	SOIL	X	$\frac{x}{x}$		$\frac{\lambda}{X}$	
Q12-4.5 Q12-15.0	08-29782-NX71P	SOIL	X	$\frac{x}{x}$	X	X	<u> </u>
Q12-10.0 Q12-20.5	08-29783-NX71Q	SOIL	X	$\frac{1}{x}$		- X	<del></del>
Q13-10.0	08-29786-NX71T	SOIL	X	$\frac{\lambda}{X}$	X	X	
Q13-15.0	08-29834-NX79A	SOIL	$\frac{\lambda}{X}$	$\frac{\lambda}{X}$		X	<del> </del>
Q13-10.0	08-29835-NX79B	SOIL	X	X	Χ	X	<del> </del>
Q14-2.5	08-29836-NX79C	SOIL	X	X		X	
Q14-2.5 Q14-6.5	08-29837-NX79D	SOIL	X	X		X	<del> </del>
Q14-0.5 Q14-15.5	08-29838-NX79E	SOIL	X	X	Х	X	
Q14-10.0 Q14-22.0	08-29839-NX79F	SOIL	X	X	$\frac{\lambda}{X}$	X	
Q14-22.0 Q15-4.0	08-29840-NX79G	SOIL	X	$\frac{\hat{x}}{x}$	^	- ^ X	<del></del>
Q15-4.0 Q15-16.0	08-29841-NX79H	SOIL	X	X	Х	<u>X</u>	
Q15-16.0 Q15-20.0	08-29842-NX79I	SOIL	X	$\frac{1}{x}$		^	
Q15-20.0 Q16-3.5	08-29844-NX79K	SOIL	X	$\frac{\hat{x}}{x}$		^_	
Q16-3.5 Q16-20.0	08-29846-NX79M	SOIL	X	X	Х	^X	
Q16-20.0 Q16-28.0	08-29847-NX79N	SOIL	X	$\frac{1}{x}$	^	$\frac{\hat{x}}{\hat{x}}$	
Q17-11.0	08-29849-NX79P	SOIL	X	X	X	- ^	
Q17-11.0 Q17-18.0	08-29850-NX79Q	SOIL	X	X		X	
		SOIL	X	X		X	
Q17-25.0	08-29851-NX79R		X	X			
SD-3	08-29852-NX79S	SOIL			X	X	
SD-4	08-29853-NX79T	SOIL	X	X		X	
SD-5	08-29854-NX79U	SOIL	Х	Х	Χ	Χ	L

# SAMPLE INDEX POS BNSF ROW Quendall Terminal - Borings

	1				TPH-G		
Sample ID	Lab Sample ID	Matrix	PAH	TPH-Dx	BTEX	Metals	Turbidity
QRM-1-1.0	08-29948-NY02A	SOIL	Х	X		Х	
QRM-1-2.0	08-29949-NY02B	SOIL	Х	Х		Х	
QRM-2-1.0	08-29950-NY02C	SOIL	X	X		X	
QRM-2-2.0	08-29951-NY02D	SOIL	Х	X		X	
QRM-3-1.0	08-29952-NY02E	SOIL	Х	X		X	
QRM-3-2.0	08-29953-NY02F	SOIL	Х	X		X	
QRM-4-1.0	08-29954-NY02G	SOIL	X	X		X	
QRM-4-2.0	08-29955-NY02H	SOIL	Х	Х		X	
QRM-5-1.0	08-29956-NY02I	SOIL	X	X		X	
QRM-5-2.0	08-29957-NY02J	SOIL	X	Х		Х	
Q4-W	08-29958-NY02K	WATER	Х	Х	Χ		Х
Q9-W	08-29959-NY02L	WATER	Х	Х	X		
Q12-W	08-29960-NY02M	WATER	X	X	X		X
Q14-W	08-29961-NY02N	WATER	Х	X	X		X
Q17-W	08-29962-NY02O	WATER	Χ	X	Χ		Χ
WD-2	08-29963-NY02P	WATER	Χ	X	X		X
TB	08-29964-NY02Q	WATER			Χ		
TB	08-30036-NY15A	WATER			Χ		
Q4-W	08-34714-OF61A	WATER				X	
Q9-W	08-34715-OF61B	WATER				Х	
Q12-W	08-34716-OF61C	WATER				X	
Q14-W	08-34717-OF61D	WATER				Х	
Q17-W	08-34718-OF61E	WATER				X	
WD-2	08-34719-OF61F	WATER				Х	

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal Borings Polycyclic Aromatic Hydrocarbon Compounds by Method SW8270D-GC/MS

This report documents the review of analytical data from the analyses of groundwater, soil, and sediment samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples		
NX66	16 Soil		
NX71	15 Soil		
NX79	18 Soil		
NY02	6 Groundwater, 10 soil		

#### I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

### II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%).

### III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are listed below.

- 2 Holding Times and Sample Preservation GC/MS Instrument Performance Check Initial Calibration (ICAL)
  - Initial Calibration (ICAL)
  - Continuing Calibration (CCAL)
    Laboratory Blanks
  - Field Blanks
- 1 Surrogate Compounds
- 2 Laboratory Control Samples (LCS/LCSD)

- 1 Matrix Spikes/Matrix Spike Duplicates (MS/MSD)
- 2 Field Replicates Internal Standards Target Analyte List Reporting Limits
- 2 Compound Identification and Reported Results
- 1 Calculation Verification

<sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Holding Times and Sample Preservation**

The validation guidance documents state that the cooler temperatures should be within an advisory temperature range of 2° to 6°C. The laboratory received the majority of the sample coolers with temperatures outside the advisory control limits of 2° to 6°C, ranging from 1.6° to 3.6°C. These temperature outliers did not impact data quality and no action was taken.

**SDG NY02:** Samples Q4-W, Q12-W, Q14-W, Q17-W, and WD-2 were re-extracted after the holding time had expired. All results for these samples were estimated (J/UJ-1).

# **Surrogates**

The percent recovery (%R) values for the surrogates were within the specified control limits for samples analyzed at dilutions of less than 10x. Several samples required dilutions greater than 10x. Surrogate recovery control limits do not apply in these cases.

# Matrix Spike/Matrix Spike Duplicates

Matrix spike/matrix spike duplicate (MS/MSD) analyses were performed at the required frequency. The MS/MSD recovery values were within the laboratory control limits, with the exception noted below. All MS/MSD relative percent difference (RPD) values were within the specified control limits.

**SDG** NX71: For QC Sample Q11-18.0, the MS %R value for fluorene was greater than the upper control limit. No qualification was necessary as the MSD recovery was in control.

# **Laboratory Control Sample/Laboratory Control Samples**

Laboratory control sample/laboratory control sample duplicates (LCS/LCSD) analyses were performed at the required frequency. The recovery values were within the specified control limits, with the exceptions noted below.

**SDG NY02:** There were no positive results for any compounds in the LCS/LCSD from the original extraction of the water samples. The water samples were re-extracted and re-analyzed, with the exception of Q9-W which had insufficient sample remaining to re-extract. The fact that no compounds were detected in the LCS/LCSD and the acceptable surrogate recoveries for the LCS/LCSD samples indicate that the spiking solution was not added. Therefore, all results for Sample Q9-W were estimated (J-10) instead of being rejected.

## **Field Replicates**

The following acceptance criteria were used to evaluate field precision: the RPD control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL (soils) or the RL (waters).

SDG NX66 and NX79: One set of field replicates was submitted: Sample Q4-15.0, submitted with SDG NX66, and the field duplicate Sample SD-3, submitted with SDG NX79. The RPD

for dibenzofuran (52.4%) exceeded the control limit. The sample and field duplicate were estimated (J-9).

**SDG NX79:** Two sets of field replicates, Samples Q13-22.0 and SD-4, and Samples Q15-16.0 and SD-5, were submitted. All field precision criteria were met.

**SDG NY02:** One set of field replicates, Samples Q12-W and WD-2, were submitted. The RPD value for naphthalene was greater than the control limit. For acenaphthene and pyrene, the difference between the sample and replicate was greater than the RL. The results for these analytes were estimated (J-9) in both samples.

# **Compound Identification and Reported Results**

Samples were screened in order to determine necessary dilutions. In some cases, one or more analytes still exceeded the instrument linear range and required additional dilution. The results for the compounds in the original analyses that were greater than the calibration range were labeled do-not-report (DNR-20). Results for all other compounds in additional dilutions were labeled do not report (DNR-11). The samples requiring additional dilutions and the associated compounds are listed below:

*SDG NX66:* Q4-2.5 (20x), Q4-15.0 (300x), Q5-18.0 (3x), Q6-4.0 (50x), Q7-4.0 (20x), Q7-5.5 (20x)

**SDG NX71:** Q9-28.0 (3x), Q10-5.0 (10x)

**SDG NX79:** Q14-6.5 (20x), Q15-4.0 (20x), SD-3 (200x)

**SDG NY02:** Q4-W (200x), Q9-W (250x)

### **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were found.

### IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, LCS/LCSD, and MS/MSD %R values, with the exceptions noted above. Precision was also acceptable as demonstrated by the field replicate, MS/MSD, and LCS/LCSD RPD values, except as previously noted.

Data were estimated based on exceeded holding times, LCS/LCSD %R outliers, and field replicate RPD outliers.

Data were labeled do-not-report (DNR) to indicate which result, from multiple dilutions and analyses, should be used. A usable result remains for all analytes in all samples, therefore completeness is unaffected.

Data labeled DNR should not be used for any purpose. All other data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal Borings Diesel and Residual Range Hydrocarbons by NWTPH-Dx

This report documents the review of analytical data from the analyses of groundwater, soil, and sediment samples and the associated field and laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NX66	16 Soil
NX71	15 Soil
NX79	18 Soil
NY02	6 Groundwater, 10 soil

#### I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

#### II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

### III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are summarized in the table below.

 Holding Times and Sample Preservation Initial Calibration (ICAL) Continuing Calibration (CCAL) Blanks (Method and Field)

2 Surrogate Compounds Laboratory Control Samples (LCS/LCSD) Matrix Spike/Matrix Spike Duplicates (MS/MSD)

Field ReplicatesReporting LimitsCompound Identification

2 Sample Results

1 Calculation Verification

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures that were outside of the recommended range of 4°C ±2°, ranging from 1.6° to 11.6°C. It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

# **Surrogate Compounds**

The surrogate recoveries for several samples could not be calculated due to necessary dilutions. No action was necessary.

**SDG NY02:** The percent recovery (%R) value for o-terphenyl was less than the lower control limit in sample Q17-W. The sample results were estimated (UJ-13) to indicate a potential low bias.

# **Field Replicates**

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

**SDG NX66 and NX79:** One set of field replicates was submitted. The parent sample, Q4-15.0, was submitted with SDG NX66. The duplicate sample, SD-3, was submitted with SDG NX79. All field precision criteria were met.

**SDG NX79:** Two sets of field replicates, samples Q13-22.0 and SD-4, and samples Q15-16.0 and SD-5, were submitted with this SDG. Field precision was acceptable.

**SDG** NY02: One set of field replicates, samples Q12-W and WD-2, were submitted with this SDG. Field precision was acceptable.

### Sample Results

**SDG NY02:** Sample Q4-W was re-extracted an re-analyzed at a 5x dilution due to a surrogate outlier in the original analysis. The re-extraction was done after the holding time had expired. The results from the original analysis should be used. The results from the re-extraction were labeled do-not-report (DNR-11).

## **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

## IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate, laboratory control sample/laboratory control sample duplicate (LCS/LCSD), and matrix spike/matrix spike duplicate (MS/MSD) recoveries, except as previously noted. Precision was also acceptable as demonstrated by the LCS/LCSD, MS/MSD, and field replicate relative percent difference (RPD) values, except as noted above.

Data were estimated because of surrogate recovery and field replicate RPD outliers.

Data were labeled as do-not-report (DNR) to indicate which results from multiple analyses should be used. A usable result remains for all analytes and all samples, therefore completeness is unaffected.

Data labeled DNR should not be used for any purpose. All other data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal Borings Gasoline Range Organics by NWTPH-Gx

This report documents the review of analytical data from the analyses of groundwater, soil, and sediment samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples			
NX66	8 Soil			
NX71	7 Soil			
NX79	9 Soil			
NY02	6 Groundwater, 1 Trip Blank			
NY15	1 Trip Blank			

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

## II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

#### III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are summarized in the table below.

 Holding Times and Sample Preservation Instrument Performance Check Initial Calibration (ICAL)
 Continuing Calibration (CCAL)

Blanks (Method and Field)
 Surrogate Compounds

Laboratory Control Samples Laboratory Duplicates

- 1 Field Replicates
- Reported Results
   Compound Identification
- 1 Calculation Verification

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $1.8^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

# Blanks (Method and Field)

**SDG NY02:** One trip blank was submitted with this SDG. There were no positive results for this blank.

**SDG** NY15: One trip blank was submitted with this SDG. There were no positive results for this blank.

# Field Replicates

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

**SDG** NX66 and NX79: One set of field replicates was submitted: Sample Q4-15.0, submitted with SDG NX66 and the field duplicate Sample SD-3, submitted with SDG NX79. Both samples were analyzed on the same instrument on the same day. All field precision criteria were met.

**SDG NX79:** Two sets of field replicates, Samples Q13-22.0 and SD-4, and Samples Q15-16.0 and SD-5, were submitted with this SDG. Field precision was acceptable.

**SDG** NY02: One set of field replicates, Samples Q12-W and WD-2 was submitted with this SDG. Field precision was acceptable.

### Reported Results

**SDG** NX71: The sample amount for each of the samples was reported incorrectly by the laboratory on the Form 1 as "mg-as-rec". The units were corrected by the validator to read "mg-dry-wt".

#### Calculation Verification

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate and laboratory control sample percent recovery values. Precision was acceptable as demonstrated by the RPD values for the laboratory and field replicate analyses.

No data were qualified for any reason.

All data, as reported, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal Borings BTEX Compounds – by SW8021Mod

This report documents the review of analytical data from the analyses of groundwater, soil, and sediment samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	G Number of Samples			
NX66	8 Soil			
NX71	7 Soil			
NX79	9 Soil			
NY02	6 Groundwater, 1 Trip Blank			
NY15	1 Trip Blank			

# I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

#### II. EDD TO HARDCOPY VERIFICATION

A complete (100%) verification of the electronic data deliverable (EDD) results was performed by comparison to the hardcopy laboratory data package. Laboratory QC results were also verified (10%). No errors were found.

## III. TECHNICAL DATA VALIDATION

The QC requirements that were reviewed are summarized in the table below.

 Holding Times and Sample Preservation Instrument Performance Check Initial Calibration (ICAL) Continuing Calibration (CCAL)

Blanks (Method and Field)
 Surrogate Compounds

Laboratory Control Samples (LCS/LCSD)

- 2 Field Replicates
- Reported Results
   Compound Identification
- 1 Calculation Verification

<sup>&</sup>lt;sup>1</sup> Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below,

# **Holding Times and Sample Preservation**

Several coolers were received at temperatures outside of the recommended range of  $4^{\circ}\text{C}$   $\pm 2^{\circ}$ , ranging from 1.6° to 11.6°C. It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

# Blanks (Method and Field)

**SDG** NY02: One trip blank was submitted with this SDG. No target analytes were detected in this blank.

**SDG** NY15: One trip blank was submitted with this SDG. No target analytes were detected in this blank.

# Field Replicates

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

**SDG** NM40: One set of field replicates, Samples Q2-D-W and WD-1 was submitted with this SDG. All field precision criteria were met.

**SDG** NM43: One set of field replicates, Samples Q1-D-23.0 and SD-2 was submitted with this SDG. All field precision criteria were met.

**SDG** NX66 and NX79: One set of field replicates was submitted: Sample Q4-15.0, submitted with SDG NX66 and the field duplicate Sample SD-3, submitted with SDG NX79. The RPD for benzene (75.2%) exceeded the control limit. The benzene results for these two samples were estimated (J-9).

**SDG** NX79: Two sets of field replicates, Samples Q13-22.0 and SD-4, and Samples Q15-16.0 and SD-5, were submitted with this SDG. Field precision was acceptable.

**SDG** NY02: One set of field replicates, Samples Q12-W and WD-2 was submitted with this SDG. Field precision was acceptable.

### **Reported Results**

**SDG NX71:** The sample amount for each of the samples was reported incorrectly by the laboratory on the Form 1 as "mg-as-rec". The units were corrected by the validator to read "mg-dry-wt".

# **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

# IV. OVERALL ASSESSMENT

As was determined by this evaluation, the laboratory followed the specified analytical method. Accuracy was acceptable, as demonstrated by the surrogate and laboratory control sample/laboratory control sample duplicate (LCS/LCSD) percent recovery values. Precision was acceptable as demonstrated by the RPD values for the LCS/LCSD and field replicate analyses, with the exception noted above.

Data were estimated based on a field precision outlier.

All data, as qualified, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal Borings Metals by Methods SW6010B, EPA 200.8, and SW7471A

This report documents the review of analytical data from the analyses of groundwater, soil, and sediment samples and the associated laboratory and field quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All Data received a full (Level IV) validation. See the Sample Index for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NX66	16 Soil
NX71	15 Soil
NX79	18 Soil
NY02	10 Soil
OF61	5 Groundwater

#### I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables, with the exception noted below. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

SDG NY02: The water samples in this SDG were re-analyzed and reported in SDG OF61.

**SDG OF61:** This SDG consists of water samples originally submitted in SDG NY02. The samples were re-analyzed by ICP-MS in order to achieve the project required detection limits. The laboratory analyzed sample Q9-W for arsenic and lead, even though these analyses were not requested on the Chain of Custody form.

#### II. EDD TO HARDCOPY VERIFICATION

A verification of the electronic data deliverables (EDD) results was performed by comparison to the hardcopy laboratory data package. No errors were found.

#### III. TECHNICAL DATA VALIDATION

The QC requirements for review are listed below.

1 Holding Times and Sample Preservation

Initial Calibration

Calibration Verification

**CRDL Standards** 

Laboratory Blanks

Laboratory Control Samples (LCS)

Matrix Spikes (MS)

**Laboratory Duplicates** 

Field Blanks

1 Field Replicates

Interference Check Samples

Serial Dilutions

ICPMS Internal Standards

1 Reported Results

1 Calculation Verification

#### **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $1.8^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

#### Field Replicates

The following acceptance criteria were used to evaluate field precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL (soils) or the RL (waters).

**SDG** NX66 and NX79: One set of field replicates was submitted: Sample Q4-15.0, submitted with SDG NX66 and the field duplicate Sample SD-3, submitted with SDG NX79. There were no positive results in either sample. Field precision was acceptable.

**SDG NX79:** Two sets of field replicates, Samples Q13-22.0 and SD-4, and Samples Q15-16.0 and SD-5, were submitted with this SDG. Field precision was acceptable.

*SDG OF61:* One set of field replicates, Samples Q12-W and WD-2, was submitted with this SDG. All field precision criteria were met.

#### **Reported Results**

Several reporting limits slightly exceeded the QAPP specified levels due to the sample size prepared, the percent total solids of the sample, and the dilution required for analysis.

**SDG NY02:** The water samples in this SDG were analyzed for arsenic and lead by Method SW6010B (ICP). The detection limits by this method do not meet the project requirements. All

Quality control results are discussed below, but no data were qualified.

<sup>&</sup>lt;sup>2</sup> Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

water samples were re-analyzed by Method 200.8 (ICP-MS) and reported in SDG OF61. The required detection limits were attained and no further action was necessary.

#### **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

#### IV. OVERALL ASSESSMENT

As determined by this evaluation, the laboratory followed the specified analytical methods. The laboratory and field duplicate RPD values indicated acceptable precision, except as noted above. Accuracy was also acceptable, as demonstrated by the matrix spike and laboratory control sample recoveries.

No data were qualified for any reason. All data, as reported, are acceptable for use.

# DATA VALIDATION REPORT Pinnacle GeoSciences POS BNSF ROW - Quendall Terminal Borings Total Solids by 160.3 and Turbidity by 180.1

This report documents the review of analytical data from the analyses of groundwater and soil samples and the associated laboratory quality control (QC) samples. Analytical Resources, Inc., Tukwila, Washington, analyzed the samples. All data received a full (Level IV) validation. See the **Sample Index** for a complete list of samples for which data were reviewed.

SDG	Number of Samples
NX66	16 Soil
NX71	15 Soil
NX79	18 Soil
NY02	5 Groundwater, 10 soil

#### I. DATA PACKAGE COMPLETENESS

The laboratory submitted all required deliverables. The laboratory followed adequate corrective action processes and all anomalies were discussed in the case narrative.

#### II. EDD TO HARDCOPY VERIFICATION

A verification of the electronic data deliverables (EDD) results was performed by comparison to the hardcopy laboratory data package. Errors in the units for the turbidity duplicate analysis for samples Q2-D-W and Q4-W were found. These were corrected in the EDD and no further action was taken.

#### III. TECHNICAL DATA VALIDATION

The QC requirements for review are listed below.

Holding Times and Sample Preservation
 Laboratory Blanks
 Laboratory Control Samples
 Matrix Spike (MS)

Laboratory Replicates

Field Replicates
 Reporting Limits

1 Calculation Verification

#### **Holding Times and Sample Preservation**

Several coolers were received at temperatures greater than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $6.8^{\circ}$  to  $11.6^{\circ}\text{C}$ . Several coolers were received at temperatures less than the recommended range of  $4^{\circ}\text{C} \pm 2^{\circ}$ , ranging from  $1.6^{\circ}$  to  $1.8^{\circ}\text{C}$ . It was determined that these temperature outliers did not impact data quality and no qualifiers were assigned.

Quality control results are discussed below, but no data were qualified.

Quality control outliers that impact the reported data were noted. Data qualifiers were issued as discussed below.

#### Field Replicates

The following acceptance criteria were used to evaluate precision: the relative percent difference (RPD) control limit is 35% for water samples and 50% for soil sample results greater than five times the reporting limit (RL). For results less than five times the RL, the difference between the sample and replicate must be less than two times the RL.

No data were qualified based on field replicate precision outliers. Users of the data should consider the impact of field precision outliers on the reported results.

**SDG NX66 and NX79:** One set of field replicates was submitted: Sample Q4-15.0, submitted with SDG NX66 and the field duplicate Sample SD-3, submitted with SDG NX79. Field precision was acceptable.

**SDG NX79:** Two sets of field replicates, Samples Q13-22.0 and SD-4, and Samples Q15-16.0 and SD-5, were submitted with this SDG. Field precision was acceptable.

**SDG NY02:** One set of field replicates, Samples Q12-W and WD-2, was submitted with this SDG. All field precision criteria were met.

#### **Calculation Verification**

Several results were verified by recalculation from the raw data. No calculation or transcription errors were noted.

#### IV. OVERALL ASSESSMENT

As determined by this evaluation, the laboratory followed the specified analytical methods. The laboratory and field replicate RPD values indicated acceptable precision. Accuracy was also acceptable, as demonstrated by the laboratory control sample recoveries.

No data were qualified for any reason. All data, as reported, are acceptable for use.



# APPENDIX A DATA QUALIFIER DEFINITIONS REASON CODES AND CRITERIA TABLES

### DATA VALIDATION QUALIFIER CODES National Functional Guidelines

The following definitions provide brief explanations of the qualifiers assigned to results in the data review process.

U	The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
J	The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.
N	The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification".
NJ	The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents the approximate concentration.
UJ	The analyte was not detected above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.
R	The sample results are rejected due to serious deficiencies in the ability to analyze the sample and meet quality control criteria. The presence or absence of the analyte cannot be verified.

The following is an EcoChem qualifier that may also be assigned during the data review process:

DNR Do not report; a more appropriate result is reported from another analysis or dilution.

### **DATA QUALIFIER REASON CODES**

1	Holding Time/Sample Preservation	
2	Chromatographic pattern in sample does not match pattern of calibration standard.	
3	Compound Confirmation	
4	Tentatively Identified Compound (TIC) (associated with NJ only)	,
5A	Calibration (initial)	
5B	Calibration (continuing)	
6	Field Blank Contamination	
7	Lab Blank Contamination (e.g., method blank, instrument, etc.)	
8	Matrix Spike(MS & MSD) Recoveries	
9	Precision (all replicates)	
10	Laboratory Control Sample Recoveries	
11	A more appropriate result is reported (associated with "R" and "DNR" only)	
12	Reference Material	
13	Surrogate Spike Recoveries (a.k.a., labeled compounds & recovery standards)	
14	Other (define in validation report)	
15	GFAA Post Digestion Spike Recoveries	
16	ICP Serial Dilution % Difference	
17	ICP Interference Check Standard Recovery	
18	Trip Blank Contamination	
19	Internal Standard Performance (e.g., area, retention time, recovery)	
20	Linear Range Exceeded	
21	Potential False Positives	
22	Elevated Detection Limit Due to Interference (i.e., laboratory, chemical and/or matrix)	

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### EcoChem Validation Guidelines for Semivolatile Analysis by GC/MS (Based on Organic NFG 1999)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE
Cooler Temperature	4°C ±2°	J(+)/UJ(-) if greater than 6 deg. C (EcoChem PJ)	1
Holding Time	Water: 7 days from collection Soil: 14 days from collection Analysis: 40 days from extraction	\frac{Water:}{J(+)/UJ(-) if ext. > 7 and < 21 days}{J(+)/R(-) if ext > 21 days (EcoChem PJ)}{\frac{Solids/Wastes}{Solids/Wastes}}: \frac{J(+)/UJ(-) if ext. > 14 and < 42 days}{J(+)/R(-) if ext. > 42 days (EcoChem PJ)}	1
		J(+)/UJ(-) if analysis >40 days	
Tuning	DFTPP Beginning of each 12 hour period Method acceptance criteria	R(+/-) all analytes in all samples associated with the tune	5A
Initial Calibration	RRF > 0.05	(EcoChem PJ, see TM-06)  If MDL= reporting limit: $J(+)/R(-) \text{ if } RRF < 0.05$	5A
(Minimum 5 stds.)		If reporting limit > MDL: note in worksheet if RRF <0.05	
	%RSD < 30%	(EcoChem PJ, see TM-06) J(+) if %RSD > 30%	5A
	RRF > 0.05	(EcoChem PJ, see TM-06)  If MDL= reporting limit:  J(+)/R(-) if RRF < 0.05	5B
Continuing Calibration (Prior to each 12 hr.		If reporting limit > MDL; note in worksheet if RRF <0.05	
shift)	%D <25%	(EcoChem PJ, see TM-06) If > +/-90%: J+/R- If -90% to -26%: J+ (high bias) If 26% to 90%: J+/UJ- (low bias)	5B
	One per matrix per batch No results > CRQL	U(+) if sample (+) result is less than CRQL and less than appropriate 5X or 10X rule (raise sample value to CRQL)	7
Method Blank		U(+) if sample (+) result is greater than or equal to CRQL and less than appropriate 5X and 10X rule (at reported sample value)	7
	No TICs present	R(+) TICs using 10X rule	7
Field Blanks (Not Required)	No results > CRQL	Apply 5X/10X rule; U(+) < action level	6

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### EcoChem Validation Guidelines for Semivolatile Analysis by GC/MS (Based on Organic NFG 1999)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON
MS/MSD (recovery)	One per matrix per batch Use method acceptance criteria	Qualify parent only unless other QC indicates systematic problems: J(+) if both %R > UCL J(+)/UJ(-) if both %R < LCL J(+)/R(-) if both %R < 10% PJ if only one %R outlier	8
MS/MSD (RPD)	One per matrix per batch Use method acceptance criteria	J(+) in parent sample if RPD > CL	9
LCS low conc. H2O SVOA	One per lab batch Within method control limits	J(+) assoc. cmpd if > UCL $J(+)/R(-)$ assoc. cmpd if < LCL $J(+)/R(-)$ all cmpds if half are < LCL	10
LCS regular SVOA (H2O & solid)	One per lab batch Lab or method control limits	J(+) if %R > UCL $J(+)/UJ(-)$ if %R <lcl <math="">J(+)/R(-) if %R &lt; 10% (EcoChem PJ)</lcl>	10
LCS/LCSD (if required)	One set per matrix and batch of 20 samples RPD < 35%	J(+)/UJ(-) assoc. cmpd. in all samples	9
Surrogates	Minimum of 3 acid and 3 base/neutral compounds Use method acceptance criteria	Do not qualify if only 1 acid and/or 1 B/N surrogate is out unless <10% J(+) if %R > UCL J(+)/UJ(-) if %R < LCL J(+)/R(-) if %R < 10%	13
Internal Standards	Added to all samples Acceptable Range: IS area 50% to 200% of CCAL area RT within 30 seconds of CC RT	J(+) if $> 200\%$ J(+)/UJ(-) if $< 50\%$ J(+)/R(-) if $< 25\%$ RT>30 seconds, narrate and Notify PM	19
Field Duplicates	Use QAPP limits. If no QAPP: Solids: RPD <50% OR absolute diff. < 2X RL (for results < 5X RL)  Aqueous: RPD <35% OR absolute diff. < 1X RL (for results < 5X RL)	Narrate and qualify if required by project (EcoChem PJ)	9
TICs	Major ions (>10%) in reference must be present in sample; intensities agree within 20%; check identification	NJ the TIC unless: R(+) common laboratory contaminants See Technical Director for ID issues	4
Quantitation/ Identification	RRT within 0.06 of standard RRT Ion relative intensity within 20% of standard All ions in std. at > 10% intensity must be present in sample	See Technical Director if outliers	14 21 (false +)

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### EcoChem Validation Guidelines for Total Petroleum Hydrocarbons-Diesel & Residual Range (Based on EPA National Functional Guidelines as applied to criteria in NWTPH-Dx, June 1997, Wa DOE & Oregon DEQ)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE
Cooler Temperature & Preservation	4°C±2°C Water: HCl to pH < 2	J(+)/UJ(-) if greater than 6 deg. C	1
Holding Time	Ext. Waters: 14 days preserved 7 days unpreserved Ext. Solids: 14 Days Analysis: 40 days from extraction	J(+)/UJ(-) if hold times exceeded J(+)/R(-) if exceeded > 3X (EcoChem PJ)	1
Initial Calibration	5 calibration points (All within 15% of true value)  Linear Regression: R <sup>2</sup> ≥0.990  If used, RSD of response factors ≤20%	Narrate if fewer than 5 calibration levels or if %R >15%  J(+)/UJ(-) if R <sup>2</sup> <0.990  J(+)/UJ(-) if %RSD > 20%	5A
Mid-range Calibration Check Std.	Analyzed before and after each analysis shift & every 20 samples.  Recovery range 85% to 115%	Narrate if frequency not met.  J(+)/UJ(-) if %R < 85%  J(+) if %R >115%	5B
Method Blank	At least one per batch (≤10 samples)	U (at the RL) if sample result is < RL & < 5X blank result.	7
	No results >RL	U (at reported sample value) if sample result is ≥ RL and < 5X blank result	7
Field Blanks (if required by project)	No results > RL	Action is same as method blank for positive results remaining in the field blank after method blank qualifiers are assigned.	6
MS samples (accuracy) (if required by project)	%R within lab control limits	Qualify parent only, unless other QC indicates systematic problems.  J(+) if both %R > upper control limit (UCL)  J(+)/UJ(-) if both %R < lower control limit (LCL)  No action if parent conc. >5X the amount spiked.  Use PJ if only one %R outlier	8
Precision: MS/MSD or LCS/LCSD or sample/dup	At least one set per batch (≤10 samples) RPD ≤ lab control limit	J(+) if RPD > lab control limits	9
LCS (not required by method)	%R within lab control limits	J(+)/UJ(-) if %R < LCL J(+) if %R > UCL J(+)/R(-) if any %R <10% (EcoChem PJ)	10

Table No.: NWTPH-Dx Revision No.: 2 Last Rev. Date: 8/13/07

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### EcoChem Validation Guidelines for Total Petroleum Hydrocarbons-Diesel & Residual Range (Based on EPA National Functional Guidelines as applied to criteria in NWTPH-Dx, June 1997, Wa DOE & Oregon DEQ)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE
Surrogates	2-fluorobiphenyl, p-terphenyl, o-terphenyl, and/or pentacosane added to all samples (inc. QC samples).  %R = 50-150%	J(+)/UJ(-) if %R < LCL J(+) if %R > UCL J(+)/R(-) if any %R <10% No action if 2 or more surrogates are used, and only one is outside control limits. (EcoChem PJ)	13
Pattern Identification	Compare sample chromatogram to standard chromatogram to ensure range and pattern are reasonable match.  Laboratory may flag results which have poor match.	J(+)	2
Field Duplicates	Use project control limits, if stated in QAPP  EcoChem default:  water: RPD < 35%  solids: RPD < 50%	Narrate (Use Professional Judgement to qualify)	9
Two analyses for one sample (dilution)	Report only one result per analyte	"DNR" (or client requested qualifier) all results that should not be reported. (See TM-04)	11

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### EcoChem Validation Guidelines for Total Petroleum Hydrocarbons-Gasoline Range (Based on EPA National Functional Guidelines as applied to criteria in NWTPH-Gx, June 1997, Wa DOE & Oregon DEQ)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE
Cooler Temperature & Preservation	4°C±2°C Water: HCl to pH < 2	J(+)/UJ(-) if greater than 6 deg. C	1
Holding Time	Waters: 14 days preserved 7 days unpreserved Solids: 14 Days	J(+)/UJ(-) if hold times exceeded J(+)/R(-) if exceeded > 3X (EcoChem PJ)	1
Initial Calibration	5 calibration points (All within 15% of true value)  Linear Regression: R <sup>2</sup> ≥0.990 If used, RSD of response factors ≤20%	Narrate if fewer than 5 calibration levels or if %R >15% J(+)/UJ(-) if R <sup>2</sup> <0.990 J(+)/UJ(-) if %RSD > 20%	5A
Mid-range Calibration Check Std.	Analyzed before and after each analysis shift & every 20 samples.  Recovery range 80% to 120%	Narrate if frequency not met.  J(+)/UJ(-) if %R < 80%  J(+) if %R >120%	5B
Method Blank	At least one per batch (≤10 samples)	U (at the RL) if sample result is < RL & < 5X blank result.	7
Wethod Dialik	No results >RL	U (at reported sample value) if sample result is ≥ RL and < 5X blank result	7
Trip Blank (if required by project)	No results >RL	Action is same as method blank for positive results remaining in trip blank after method blank qualifiers are assigned.	18
Field Blanks (if required by project)	No results > RL	Action is same as method blank for positive results remaining in field blank after method <b>and</b> trip blank qualifiers are assigned.	6
MS samples (accuracy) (if required by project)	%R within lab control limits	Qualify parent only, unless other QC indicates systematic problems.  J(+) if both %R > upper control limit (UCL)  J(+)/UJ(-) if both %R < lower control limit (LCL)  No action if parent conc. >5X the amount spiked.  Use PJ if only one %R outlier	8
Precision: MS/MSD or LCS/LCSD or sample/dup	At least one set per batch (≤10 samples) RPD ≤ lab control limit	J(+) if RPD > lab control limits	9

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### EcoChem Validation Guidelines for Total Petroleum Hydrocarbons-Gasoline Range (Based on EPA National Functional Guidelines as applied to criteria in NWTPH-Gx, June 1997, Wa DOE & Oregon DEQ)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE
LCS (not required by method)	%R within lab control limits	J(+)/UJ(-) if %R < LCL J(+) if %R > UCL J(+)/R(-) if any %R <10% (EcoChem PJ)	10
Surrogates	Bromofluorobenzene and/or 1,4-difluorobenzene added to all samples (inc. QC samples).  %R = 50-150%	J(+)/UJ(-) if %R < LCL J(+) if %R >UCL J(+)/R(-) if any %R <10% No action if 2 or more surrogates are used, and only one is outside control limits. (EcoChem PJ)	13
Pattern Identification	Compare sample chromatogram to standard chromatogram to ensure range and pattern are reasonable match.  Laboratory may flag results which have poor match.	J(+)	2
Field Duplicates	Use project control limits, if stated in QAPP  EcoChem default: water: RPD < 35% solids: RPD < 50%	Narrate outliers  If required by project, qualify with J(+)/UJ(-)	9
Two analyses for one sample (e.g., dilution)	Report only one result per analyte	"DNR" (or client requested qualifier) all results that should not be reported. (See TM-04)	11

Table No.: NFG-Pest PCB Revision No.: 4 Last Rev. Date: 8/23/07

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### EcoChem Validation Guidelines for Pesticides/PCBs by GC/ECD (Based on Organic NFG 1999 & EPA SW-846 Method 8081/8082)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON
Cooler Temperature	4°C ±2°	J(+)/UJ(-) if greater than 6 deg. C (EcoChem PJ)	1
Holding Time	Water: 7 days from collection Soil: 14 days from collection Analysis: 40 days from extraction	J(+)/UJ(-) if ext/analyzed > HT J(+)/R(-) if ext/analyzed > 3X HT (EcoChem PJ)	1
Resolution Check	Beginning of ICAL Sequence Within RTW Resolution >90%	Narrate (Use Professional Judgement to qualify)	14
Instrument Performance (Breakdown)	DDT Breakdown: < 20% Endrin Breakdown: <20% Combined Breakdown: <30% Compounds within RTW	J(+) DDT NJ(+) DDD and/or DDE R(-) DDT - If (+) for either DDE or DDD  J(+) Endrin NJ(+) EK and/or EA R(-) Endrin - If (+) for either EK or EA	5A
Retention Times	Surrogates: TCX (+/- 0.05); DCB (+/- 0.10) Target compounds: elute before heptachlor epoxide (+/- 0.05) elute after heptachlor epoxide (+/- 0.07)	NJ(+)/R(-) results for analytes with RT shifts For full DV, use PJ based on examination of raw data	5B
Initial Calibration	Pesticides: Low=CRQL, Mid=4X, High=16X Multiresponse - one point Calibration %RSD<20% %RSD<30% for surr; two comp. may exceed if <30% Resolution in Mix A and Mix B >90%	J(+)/UJ(-)	5A
Continuing Calibration	Alternating PEM standard and INDA/INDB standards every 12 hours (each preceeded by an inst. Blank) %D < 25%  Resolution >90% in IND mixes; 100% for PEM	J(+)/UJ(-) J(+)R(-) if %D > 90%  PJ for resolution	5B
Method Blank	One per matrix per batch No results > CRQL	U(+) if sample result is < CRQL and < 5X rule (raise sample value to CRQL)  U(+) if sample result is > or equal to CRQL and	7
Instrument Blanks	Analyzed at the beginning of every 12 hour sequence No analyte > 1/2 CRQL	< 5X rule (at reported sample value)  Same as Method Blank	7
Field Blanks	Not addressed by NFG No results > CRQL	Apply 5X rule; U(+) < action level	6

Table No.: NFG-Pest PCB Revision No.: 4 Last Rev. Date: 8/23/07 Page: 2 of 2

### EcoChem Validation Guidelines for Pesticides/PCBs by GC/ECD (Based on Organic NFG 1999 & EPA SW-846 Method 8081/8082)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON
MS/MSD (recovery)	One set per matrix per batch Method Acceptance Criteria	Qualify parent only unless other QC indicates systematic problems: J(+) if both %R > UCL J(+)/UJ(-) if both %R < LCL J(+)/R(-) if both %R < 10% PJ if only one %R outlier	8
MS/MSD (RPD)	One set per matrix per batch Method Acceptance Criteria	J(+) in parent sample if RPD > CL	9
LCS	One per SDG Method Acceptance Criteria	J(+) if %R > UCL J(+)/UJ(-) if %R < LCL J(+)/R(-) using PJ if %R < <lcl (<="" 10%)<="" td=""><td>10</td></lcl>	10
LCS/LCSD (if required)	One set per matrix and batch of 20 samples RPD < 35%	J(+)/UJ(-) assoc. cmpd. in all samples	9
Surrogates	TCX and DCB added to every sample %R = 30-150%	J(+)/UJ(-) if both %R = 10 - 60% J(+) if both >150% J(+)/R(-) if any %R <10%	13
Quantitation/ Identification	Quantitated using ICAL calibration factor (CF)  RPD between columns <40%	J(+) if RPD = 40 - 60% NJ(+) if RPD >60% EcoChem PJ - See TM-08	3
Two analyses for one sample	Report only one result per analyte	"DNR" results that should not be used to avoid reporting two results for one sample	11
Sample Clean-up	GPC required for soil samples Florisil required for all samples Sulfur is optional Clean-up standard check %R within CLP limits	J(+)/UJ(-) if %R < LCL J(+) if %R > UCL	14
Field Duplicates	Use QAPP limits. If no QAPP: Solids: RPD <50% OR absolute diff. < 2X RL (for results < 5X RL) Aqueous: RPD <35% OR absolute diff. < 1X RL (for results < 5X RL)	Narrate (Qualifiy if required by project QAPP)	9

Table No.: NFG-ICP Revision No.: draft Last Rev. Date: draft Page: 1 of 2

### EcoChem Validation Guidelines for Metals Analysis by ICP (Based on Inorganic NFG 1994 & 2004)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE	
Cooler Temperature and Preservation	THOSE DISCOVED METALS. O ASIM THESE & DESCRIVE SITES TO THE STATE TO THE STATE OF T			
Holding Time	180 days from date sampled Frozen tissues - HT extended to 2 years	J(+)/UJ(-) if holding time exceeded	1	
Initial Calibration	Blank + minimum 1 standard If more than 1 standard, r > 0.995	J(+)/UJ(-) if r < 0.995 (multi point cal)	5A	
Initial Calibration Verification (ICV)	J(+)/UJ(-) if %R 75-89% nitial Calibration Independent source analyzed immediately after calibration J(+) if %R = 111-125%			
Continuing Calibration Verification (CCV)	Every ten samples, immediately following ICV/ICB and at end of run %R within ±10% of true value	J(+)/UJ(-) if %R = 75-89% J(+) if %R 111-125% R(+) if %R > 125% R(+/-) if %R < 75%	5B	
Initial and Continuing Calibration Blank (ICB/CCB)	After each ICV and CCV every ten samples and end of run   blank   < IDL (MDL)	Action level is 5x absolute value of blank conc. For (+) blanks, U(+) results < action level For (-) blanks, J(+)/UJ(-) results < action level (Refer to TM-02 for additional information)	7	
Reporting Limit Standard	2x RL analyzed beginning of run Not required for Al, Ba, Ca, Fe, Mg, Na, K %R = 70%-130% (50%-150% Sb, Pb, Tl)	R(-)/J(+) < $2x$ RL if %R < $50\%$ (< $30\%$ Sb, Pb, Tl) J(+) < $2x$ RL, UJ(-) if %R 50-69% (30-49% Sb, Pb,Tl) J(+) < $2x$ RL if %R 130-180% (150-200% Sb, Pb, Tl) R(+) < $2x$ RL if %R > $180\%$ (200% Sb, Pb, Tl)	14	
Interference Check Samples (ICSA/ICSAB)	ICSAB %R 80 - 120% for all spiked elements   ICSA   < MDL for all unspiked elements except: K, Na	For samples with Al, Ca, Fe, or Mg > ICS levels R(+/-) if %R < 50% J(+) if %R >120% J(+)/UJ(-) if %R= 50 to 79% Use Professional Judgment for ICSA to determine if bias is present see TM-09 for additional details	17	
Method Blank	One per matrix per batch (batch not to exceed 20 samples) blank < MDL	Action level is 5x blank concentration U(+) results < action level	7	

Table No.: NFG-ICP Revision No.: draft Last Rev. Date: draft Page: 2 of 2

### EcoChem Validation Guidelines for Metals Analysis by ICP (Based on Inorganic NFG 1994 & 2004)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE
	One per matrix per batch		
Laboratory Control Sample (LCS)	Blank Spike: %R within 80-120%	R(+/-) if %R < 50% J(+)/UJ(-) if %R = 50-79% J(+) if %R >120%	10
	CRM: Result within manufacturer's certified acceptance range or project guidelines	J(+)/UJ(-) if < LCL, J(+) if > UCL	
Matrix Spikes	J(+) if %R > 125%  J(+)/UJ(-) if %R < 75%		8
Post-digestion Spike	If Matrix Spike is outside 75-125%, spike at twice the sample conc.	No qualifiers assigned based on this element	
Laboratory Duplicate (or MS/MSD)	One per matrix per batch RPD < 20% for samples > 5x RL Diff < RL for samples >RL and < 5x RL (Diff < 2x RL for solids)	J(+)/UJ(-) if RPD > 20% or diff > RL (2x RL for solids) qualify all samples in batch	9
Serial Dilution	5x dilution one per matrix %D < 10% for original sample conc. > 50x MDL	J(+)/UJ(-) if %D >10% qualify all samples in batch	16
Field Blank	Blank < MDL	Action level is 5x blank conc. U(+) sample values < action level in associated field samples only	6
Field Duplicate	For results > 5x RL: Water: RPD < 35% Solid: RPD < 50% For results < 5 x RL: Water: Diff < RL Solid: Diff < 2x RL	J(+)/UJ(-) in parent samples only	
Linear Range	Sample concentrations must fall within range	J values over range	_ 20

Table No.: NFG-ICPMS Revision No.: Draft Last Rev. Date: Draft Page: 1 of 2

### EcoChem Validation Guidelines for Metals Analysis by ICP-MS (Based on Inorganic NFG 1994 & 2004)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE			
Cooler Temperature and Preservation	Preservation For Dissolved Metals: 0.45um filter & preserve after filtration J(+)/UJ(-) if pH preservation requirements are not met					
Holding Time	180 days from date sampled Frozen tissues - HT extended to 2 years	J(+)/UJ(-) if holding time exceeded	1			
Tune	Prior to ICAL monitoring compounds analyzed 5 times wih Std Dev. ≤ 5% Tune Tune Resolution < 0.9 AMU @ 10% peak height or <0.75 amu @ 5% peak height  Prior to ICAL Use Professional Judgment to evaluate tune J(+)/UJ(-) if tune criteria not met					
Initial Calibration	Blank + minimum 1 standard If more than 1 standard, r>0.995	J(+)/UJ(-) if r<0.995 (for multi point cal)	5A			
Initial Calibration Verification (ICV)	Independent source analyzed immediately after calibration  ## Within +10% of true value    J(+)/UJ(-) if %R 75-89%   J(+) if %R = 111-125%					
Continuing Calibration Verification (CCV)	Every ten samples, immediately following ICV/ICB and at end of run ±10% of true value	J(+)/UJ(-) if %R = 75-89% J(+) if %R 111-125% R(+) if %R > 125% R(+/-) if %R < 75%				
Initial and Continuing Calibration Blanks (ICB/CCB)	After each ICV and CCV every ten samples and end of run   blank   < IDL (MDL)	Action level is 5x absolute value of blank conc.  For (+) blanks, U(+) results < action level  For (-) blanks, J(+)/UJ(-) results < action level  refer to <b>TM-02</b> for additional details	7			
Reporting Limit Standard (CRI)	2x RL analyzed beginning of run Not required for Al, Ba, Ca, Fe, Mg, Na, K %R = 70%-130% (50%-150% Co,Mn, Zn)	R(-),(+) < 2x RL if %R < 50% (< 30% Co,Mn, Zn) J(+) < 2x RL, UJ(-) if %R 50-69% (30%-49% Co,Mn, Zn) J(+) < 2x RL if %R 130%-180% (150%-200% Co,Mn, Zn) R(+) < 2x RL if %R > 180% (200% Co, Mn, Zn)	14			
Interference Check Samples (ICSA/ICSAB) Required by SW 6020, but not 200.8 ICSAB %R 80% - 120% for all spiked elements J( MDL) for all unspiked elements Use Profess		For samples with AI, Ca, Fe, or Mg > ICS levels R(+/-) if %R < 50% J(+) if %R >120% J(+)/UJ(-) if %R = 50% to 79% Use Professional Judgment for ICSA to determine if bias is present see TM-09 for additional details	17			
Method Blank	One per matrix per batch (batch not to exceed 20 samples) blank < MDL	Action level is 5x blank concentration U(+) results < action level	7			
Laboratory Control	One per matrix per batch Blank Spike: %R within 80%-120%	R(+/-) if %R < 50% J(+)/UJ(-) if %R = 50-79% J(+) if %R >120%	10			
Sample (LCS)	CRM: Result within manufacturer's certified acceptance range or project guidelines	J(+)/UJ(-) if < LCL, J(+) if > UCL	-			

Table No.: NFG-ICPMS Revision No.: Draft Last Rev. Date: Draft Page: 2 of 2

### EcoChem Validation Guidelines for Metals Analysis by ICP-MS (Based on Inorganic NFG 1994 & 2004)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE
Matrix Spike/ Matrix Spike Duplicate (MS/MSD)	Duplicate 75-125% for samples where results J(+)/R(-) if %R<30% or		
Post-digestion Spike	If Matrix Spike is outside 75-125%, Spike parent sample at 2x the sample conc.	No qualifiers assigned based on this element	
Laboratory Duplicate (or MS/MSD)	One per matrix per batch RPD < 20% for samples > 5x RL Diff < RL for samples > RL and < 5 x RL (Diff < 2x RL for solids)	J(+)/UJ(-) if RPD > 20% or diff > RL all samples in batch	9
Serial Dilution	5x dilution one per matrix %D < 10% for original sample values > 50x MDL	J(+)/UJ(-) if %D >10% All samples in batch	16
Internal Standards	Every sample SW6020: 60%-125% of cal blank IS 200.8: 30%-120% of cal blank IS	J (+)/UJ (-) all analytes associated with IS outlier	19
Field Blank	Blank < MDL	Action level is 5x blank conc. U(+) sample values < AL in associated field samples only	6
Field Duplicate	For results > 5x RL: Water: RPD < 35% Solid: RPD < 50% For results < 5 x RL: Water: Diff < RL Solid: Diff < 2x RL	J(+)/UJ(-) in parent samples only	9
Linear Range	Sample concentrations must fall within range	J values over range	20

Table No.: NFG-HG Revision No.: draft Last Rev. Date: draft Page: 1 of 2

### EcoChem Validation Guidelines for Mercury Analysis by CVAA (Based on Inorganic NFG 1994 & 2004)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE		
Cooler Temperature and Preservation					
Holding Time	28 days from date sampled Frozen tissues: HT extended to 6 months  J(+)/UJ(-) if holding time exceeded				
Initial Calibration	Blank + 4 standards, one at RL r > 0.995	J(+)/UJ(-) if r<0.995	5A		
Initial Calibration Verification (ICV)	Independent source analyzed immediately after    J(+)/UJ(-) if %R = 65%-79%				
Continuing Calibration Verification (CCV)	ntinuing Calibration  Every ten samples, immediately following  J(+)/UJ(-) if %R = 65%-79%  ICV/ICB and at end of run  I/+) if %R = 121-135%				
Initial and Continuing Calibration Blanks (ICB/CCB)	Calibration Blanks every ten samples and end of run  For (+) blanks, U(+) results < action level				
Reporting Limit Standard (CRA)	Reporting Limit Standard  R(-),(+)<2xRL if %R <50%  J(+)<2x RL, UJ(-) if %R 50-69%  J(+)<2x RL, UJ(-) if %R 130-180%				
Method Blank	One per matrix per batch (batch not to exceed 20 samples) blank < MDL	Action level is 5x blank concentration U(+) results < action level	7		
	One per matrix per batch				
Laboratory Control Sample (LCS)	Blank Spike: %R within 80-120%	R(+/-) if %R < 50% J(+)/UJ(-) if %R = 50-79% J(+) if %R >120%	10		
	CRM: Result within manufacturer's certified acceptance range or project guidelines	J(+)/UJ(-) if < LCL, J(+) if > UCL			
Matrix Spike/Matrix Spike Duplicate (MS/MSD)	One per matrix per batch 5% frequency 75-125% for samples less than 4x spike level	J(+) if %R>125% J(+)/UJ(-) if %R <75% J(+)/R(-) if %R<30% all samples in batch	8		
Laboratory Duplicate (or MS/MSD)	One per matrix per batch atory Duplicate RPD < 20% for samples > 5x RL J(+)/UJ(-) if RPD > 20% or diff > RL				

Table No.: NFG-HG Revision No.: draft Last Rev. Date: draft Page: 2 of 2

### EcoChem Validation Guidelines for Mercury Analysis by CVAA (Based on Inorganic NFG 1994 & 2004)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE
Field Blank Blank < MDL		Action level is 5x blank conc. U(+) sample values < action level in associated field samples only	6
Field Duplicate	For results > 5x RL: Water: RPD < 35% Solid: RPD < 50% For results < 5x RL: Water: Diff <rl 2x="" <="" diff="" rl<="" solid:="" td=""><td>J(+)/UJ(-) in parent samples only</td><td>9</td></rl>	J(+)/UJ(-) in parent samples only	9

Table No.: Eco-Conv Revision No.: 0

Last Rev. Date: FINAL DRAFT

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### EcoChem Validation Guidelines for Conventional Chemistry Analysis (Based on EPA Standard Methods)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE		
Cooler Temperature and Preservation	Preservation Preservation: Method Specific quality for coole temp outliers  J(+)/UJ(-) if preservation requirements not met				
Holding Time	Holding Time  Method Specific  Method Specific  J(+)/UJ(-) if holding time exceeded J(+)/R(-) if HT exceeded by > 3X				
Initial Calibration	Method specific r>0.995	Use professional judgment J(+)/UJ(-) for r < 0.995	5A		
Initial Calibration Verification (ICV)	Where applicable to method Independent source analyzed immediately after calibration %R method specific, usually 90% - 110%	R(+/-) if %R significantly < LCL J(+)/UJ(-) if %R < LCL J(+) if %R > UCL R(+) if %R significantly > UCL	5A		
Continuing Cal Verification (CCV)					
Initial and Continuing Cal Blanks (ICB/CCB)	Where applicable to method Action level is 5x absolute value of blank Initial and Continuing After each ICV and CCV every ten For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks, U(+) results < action level is 5x absolute value of blank For (+) blanks < action level is 5x absolute value of blank For (+) bl		7		
Method Blank	One per matrix per batch (not to exceed 20 samples) blank < MDL  Action level is 5x absolute value of blank conc. For (+) blk value, U(+) results < action level For (-) blk value, J(+)/UJ(-) results < action level		7		
Laboratory Control	Waters: One per matrix per batch %R (80-120%)	R(+/-) if %R < 50% J(+)/UJ(-) if %R = 50-79% J(+) if %R >120%	10		
Sample	Soils: One per matrix per batch Result within manufacturer's certified acceptance range	J(+)/UJ(-) if < LCL, J(+) if > UCL	10		
Matrix Spike	One per matrix per batch; 5% frequency J(+) if %R > 125% or < 75%				
Laboratory Duplicate	One per matrix per batch RPD <20% for samples > 5x RL Diff <rl for="" samples="">RL and &lt;5 x RL (may use RPD &lt; 35%, Diff &lt; 2X RL for solids)</rl>	J(+)/UJ(-) if RPD > 20% or diff > RL all samples in batch	9		

Table No.: Eco-Conv Revision No.: 0

Last Rev. Date: FINAL DRAFT

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### EcoChem Validation Guidelines for Conventional Chemistry Analysis (Based on EPA Standard Methods)

VALIDATION QC ELEMENT	ACCEPTANCE CRITERIA	ACTION	REASON CODE
Field Blank blank < MDL		Action level is 5x blank conc. U(+) sample values < action level in associated field samples only	6
Field Duplicate	For results > 5X RL: Water: RPD < 35% Solid: RPD < 50% For results < 5 x RL: Water: Diff <rl 2x="" <="" diff="" rl<="" solid:="" td=""><td>J(+)/UJ(-) in parent samples only</td><td>9</td></rl>	J(+)/UJ(-) in parent samples only	9



### APPENDIX B QUALIFIED DATA SUMMARY TABLE

				T		Lab	<u> </u>	
Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	DV Qual RC
Q1-A-1.0	08-21754-NM43A	SW8270D	Phenanthrene	1100000	ug/kg	E	DNR	20
Q1-A-1.0	08-21754-NM43A	SW8270D	Pyrene	1300000		E	DNR	20
Q1-A-1.0	08-21754-NM43ADL	SW8270D	1,2,4-Trichlorobenzene	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	1,2-Dichlorobenzene	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	1,3-Dichlorobenzene	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	1,4-Dichlorobenzene	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	1-Methylnaphthalene	130000			DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2,2'-Oxybis(1-Chloropropane)	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2,4,5-Trichlorophenol	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2,4,6-Trichlorophenol	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2,4-Dichlorophenol	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2,4-Dimethylphenol	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2,4-Dinitrophenol	220000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2,4-Dinitrotoluene	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2,6-Dinitrotoluene	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2-Chloronaphthalene	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2-Chlorophenol	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2-Methylnaphthalene	170000			DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2-Methylphenol	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2-Nitroaniline	110000	ug/kg	U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	2-Nitrophenol	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	3,3'-Dichlorobenzidine	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	3-Nitroaniline	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	4,6-Dinitro-2-Methylphenol	220000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	4-Bromophenyl-phenylether	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	4-Chloro-3-methylphenol	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	4-Chloroaniline	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	4-Chlorophenyl-phenylether	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL 08-21754-NM43ADL	SW8270D SW8270D	4-Methylphenol 4-Nitroaniline	22000		U	DNR	11 11
Q1-A-1.0 Q1-A-1.0	08-21754-NM43ADL	SW8270D	4-Nitrophenol	110000 110000		1 0	DNR DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Acenaphthene	150000		1-0-	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Acenaphthylene	22000		Tu	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Anthracene	220000		<del> </del> -	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Benzo(a)anthracene	520000	ug/kg	╂───	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Benzo(a)pyrene	640000		+	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Benzo(b)fluoranthene	340000		<del> </del>	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Benzo(g,h,i)perylene	230000		<del> </del>	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Benzo(k)fluoranthene	400000			DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Benzoic Acid	220000		1-0-1	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Benzyl Alcohol	22000		T Ü	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	bis(2-Chloroethoxy) Methane	22000		T U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Bis-(2-Chloroethyl) Ether	22000		1 0	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	bis(2-Ethylhexyl)phthalate	22000		Ū	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Butylbenzylphthalate	22000		Ū	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Carbazole	42000			DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Chrysene	670000		1	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Dibenz(a,h)anthracene	95000			DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Dibenzofuran	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Diethylphthalate	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Dimethylphthalate	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Di-n-Butylphthalate	22000	ug/kg	U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Di-n-Octyl phthalate	22000	ug/kg	U	DNR	11

Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Lab Flag		DV Qual RC
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Fluoranthene	630000			DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Fluorene	110000			DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Hexachlorobenzene	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Hexachlorobutadiene	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Hexachlorocyclopentadiene	110000		<u> U</u>	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Hexachloroethane	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Indeno(1,2,3-cd)pyrene	180000			DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Isophorone	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Naphthalene	250000			DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Nitrobenzene	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	N-Nitroso-Di-N-Propylamine	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	N-Nitrosodiphenylamine	22000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Pentachlorophenol	110000		U	DNR	11
Q1-A-1.0	08-21754-NM43ADL	SW8270D	Phenol	22000		U	DNR	11
Q1-B-1.0	08-21756-NM43C	SW8270D	Benzo(a)anthracene	690000		E	DNR	20
Q1-B-1.0	08-21756-NM43C	SW8270D	Benzo(a)pyrene	680000		E	DNR	20
Q1-B-1.0	08-21756-NM43C	SW8270D	Benzo(b)fluoranthene	440000		E	DNR	20
Q1-B-1.0	08-21756-NM43C	SW8270D	Benzo(k)fluoranthene	390000		E	DNR	20
Q1-B-1.0	08-21756-NM43C	SW8270D	Chrysene	680000		E	DNR	20
Q1-B-1.0	08-21756-NM43C	SW8270D	Fluoranthene	640000		E	DNR	20
Q1-B-1.0	08-21756-NM43C	SW8270D	Phenanthrene	1000000		Е	DNR	20
Q1-B-1.0	08-21756-NM43C	SW8270D	Pyrene	1400000		E	DNR	20
Q1-B-1.0	08-21756-NM43CDL	SW8270D	1,2,4-Trichlorobenzene	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	1,2-Dichlorobenzene	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	1,3-Dichlorobenzene	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	1,4-Dichlorobenzene	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	1-Methylnaphthalene	150000			DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2,2'-Oxybis(1-Chloropropane)	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2,4,5-Trichlorophenol	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2,4,6-Trichlorophenol	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2,4-Dichlorophenol	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2,4-Dimethylphenol	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2,4-Dinitrophenol	190000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2,4-Dinitrotoluene	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2,6-Dinitrotoluene	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2-Chloronaphthalene	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2-Chlorophenol	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2-Methylnaphthalene	200000		<u> </u>	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2-Methylphenol	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2-Nitroaniline	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	2-Nitrophenol	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	3,3'-Dichlorobenzidine	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	3-Nitroaniline	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	4,6-Dinitro-2-Methylphenol	190000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	4-Bromophenyl-phenylether	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	4-Chloro-3-methylphenol	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	4-Chloroaniline	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	4-Chlorophenyl-phenylether	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	4-Methylphenol	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	4-Nitroaniline	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	4-Nitrophenol	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Acenaphthene	190000		<del>                                     </del>	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Acenaphthylene	19000	ug/kg	U	DNR	11

Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Lab Flag	DV Qual	DV Qual RC
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Anthracene	280000		1 lag	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Benzo(g,h,i)perylene	260000		<del> </del> -	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Benzoic Acid	190000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Benzyl Alcohol	19000		Ü	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	bis(2-Chloroethoxy) Methane	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Bis-(2-Chloroethyl) Ether	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	bis(2-Ethylhexyl)phthalate	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Butylbenzylphthalate	19000		<del>                                     </del>	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Carbazole	47000		<del>                                     </del>	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Dibenz(a,h)anthracene	110000			DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Dibenzofuran	19000	ug/kg ug/kg	l u	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Diethylphthalate	19000	ug/kg ug/kg	T Ü	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Dimethylphthalate	19000	ug/kg	T U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Di-n-Butylphthalate	19000		T U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Di-n-Octyl phthalate	19000		T U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Fluorene	150000		<del>                                     </del>	DNR	<del></del>
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Hexachlorobenzene	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Hexachlorobutadiene	19000		Ü	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Hexachlorocyclopentadiene	96000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Hexachloroethane	19000		1 0	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Indeno(1,2,3-cd)pyrene	190000		<del>                                     </del>	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Isophorone	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Naphthalene	280000		<del>                                     </del>	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Nitrobenzene	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	N-Nitroso-Di-N-Propylamine	96000		Ü	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	N-Nitrosodiphenylamine	19000		U	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Pentachlorophenol	96000		l ü	DNR	11
Q1-B-1.0	08-21756-NM43CDL	SW8270D	Phenol	19000		l ü	DNR	11
Q1-D-1.0	08-21761-NM43H	SW8270D	Benzoic Acid		ug/kg ug/kg	U	UJ	8
Q1-D-9.0	08-21763-NM43J	SW8270D	Acenaphthene	51000		E	DNR	20
Q1-D-9.0	08-21763-NM43J	SW8270D	Fluoranthene	140000	ug/kg ug/kg	E	DNR	20
Q1-D-9.0	08-21763-NM43J	SW8270D	Phenanthrene	160000		E	DNR	20
Q1-D-9.0	08-21763-NM43J	SW8270D	Pyrene	110000		E	DNR	20
Q1-D-9.0	08-21763-NM43JDL	SW8270D	1,2,4-Trichlorobenzene		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	1,2-Dichlorobenzene		ug/kg	U	DNR	<del>-                                      </del>
Q1-D-9.0	08-21763-NM43JDL	SW8270D	1,3-Dichlorobenzene		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	1,4-Dichlorobenzene		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	1-Methylnaphthalene	13000		<del>                                     </del>	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2,2'-Oxybis(1-Chloropropane)		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2,4,5-Trichlorophenol	19000		Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2,4,6-Trichlorophenol	19000		Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2,4-Dichlorophenol	19000		Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2,4-Dimethylphenol		ug/kg	Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2,4-Dinitrophenol	37000	<u> </u>	Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2,4-Dinitrotoluene	19000		Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2,6-Dinitrotoluene	19000		Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2-Chloronaphthalene		ug/kg	Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2-Chlorophenol	3700		Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2-Methylnaphthalene	10000		┝┷┥	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2-Methylphenol	3700		U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2-Nitroaniline	19000		Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	2-Nitrophenol	3700		U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	3,3'-Dichlorobenzidine	19000		Ü	DNR	11
טיפ-ח-וא	TOO-T I TOO-ININI-1990FF	DVVOZIOD	O,O ~DIGINOTODETIZIDINE	19000	ugmg		DIAI.	11

Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Lab Flag	DV Qual	DV Qual RC
Q1-D-9.0	08-21763-NM43JDL	SW8270D	3-Nitroaniline		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	4,6-Dinitro-2-Methylphenol	37000		Ι <del>ΰ</del>	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	4-Bromophenyl-phenylether		ug/kg	<del>l ü</del>	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	4-Chloro-3-methylphenol	19000		T U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	4-Chloroaniline	19000		l Ū	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	4-Chlorophenyl-phenylether		ug/kg	T Ū	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	4-Methylphenol		ug/kg	Ū	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	4-Nitroaniline	19000		U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	4-Nitrophenol	19000		Ū	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Acenaphthylene		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Anthracene	26000			DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Benzo(a)anthracene	35000			DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Benzo(a)pyrene	27000			DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Benzo(b)fluoranthene	16000			DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Benzo(g,h,i)perylene	16000			DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Benzo(k)fluoranthene	23000			DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Benzoic Acid	37000		U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Benzyl Alcohol		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	bis(2-Chloroethoxy) Methane		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Bis-(2-Chloroethyl) Ether		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	bis(2-Ethylhexyl)phthalate		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Butylbenzylphthalate	3700	ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Carbazole	7100	ug/kg		DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Chrysene	45000			DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Dibenz(a,h)anthracene		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Dibenzofuran	22000			DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Diethylphthalate		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Dimethylphthalate	3700	ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Di-n-Butylphthalate	3700	ug/kg	Ū	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Di-n-Octyl phthalate	3700	ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Fluorene	42000	ug/kg		DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Hexachlorobenzene	3700	ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Hexachlorobutadiene	3700	ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Hexachlorocyclopentadiene	19000	ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Hexachloroethane	3700	ug/kg	Ü	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Indeno(1,2,3-cd)pyrene	13000	ug/kg		DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Isophorone		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Naphthalene	18000	<u> </u>		DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Nitrobenzene		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	N-Nitroso-Di-N-Propylamine	19000		U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	N-Nitrosodiphenylamine		ug/kg	U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Pentachlorophenol	19000		U	DNR	11
Q1-D-9.0	08-21763-NM43JDL	SW8270D	Phenol	3700		U	DNR	11
Q1-D-15.0	08-21764-NM43K	SW8270D	1-Methylnaphthalene	43000		E	DNR	20
Q1-D-15.0	08-21764-NM43K	SW8270D	2-Methylnaphthalene	62000		E	DNR	20
Q1-D-15.0	08-21764-NM43K	SW8270D	Acenaphthene	62000		Е	DNR	20
Q1-D-15.0	08-21764-NM43K	SW8270D	Anthracene	21000		E	DNR	20
Q1-D-15.0	08-21764-NM43K	SW8270D	Dibenzofuran	19000		E	DNR	20
Q1-D-15.0	08-21764-NM43K	SW8270D	Fluoranthene	78000		Е	DNR	_20
	08-21764-NM43K	SW8270D	Fluorene	41000		E	DNR	20
	08-21764-NM43K	SW8270D	Naphthalene	15000		Е	DNR	20
	08-21764-NM43K	SW8270D	Phenanthrene	170000		ES	DNR	20
Q1-D-15.0	08-21764-NM43K	SW8270D	Pyrene	56000	ug/kg	E	DNR	20

						Lab		
Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag		DV Qual RC
Q1-D-15.0	08-21764-NM43KDL	SW8270D	1,2,4-Trichlorobenzene		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	1,2-Dichlorobenzene		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	1,3-Dichlorobenzene		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	1,4-Dichlorobenzene		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2,2'-Oxybis(1-Chloropropane)		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2,4,5-Trichlorophenol	24000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2,4,6-Trichlorophenol	24000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2,4-Dichlorophenol	24000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2,4-Dimethylphenol		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2,4-Dinitrophenol	47000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2,4-Dinitrotoluene	24000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2,6-Dinitrotoluene	24000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2-Chloronaphthalene		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2-Chlorophenol		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2-Methylphenol		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2-Nitroaniline	24000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	2-Nitrophenol		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	3,3'-Dichlorobenzidine	24000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	3-Nitroaniline	24000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	4,6-Dinitro-2-Methylphenol	47000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	4-Bromophenyl-phenylether		ug/kg	U	DNR	11 11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	4-Chloro-3-methylphenol	24000		U	DNR	
Q1-D-15.0	08-21764-NM43KDL	SW8270D	4-Chloroaniline	24000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	4-Chlorophenyl-phenylether		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	4-Methylphenol		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D SW8270D	4-Nitroaniline 4-Nitrophenol	24000		U	DNR DNR	11
Q1-D-15.0 Q1-D-15.0	08-21764-NM43KDL 08-21764-NM43KDL	SW8270D	Acenaphthylene	24000	ug/kg ug/kg	U	DNR	11
Q1-D-15.0 Q1-D-15.0	08-21764-NM43KDL	SW8270D	Benzo(a)anthracene	11000		<u> </u>	DNR	11
Q1-D-15.0 Q1-D-15.0	08-21764-NM43KDL	SW8270D	Benzo(a)pyrene	4700		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Benzo(b)fluoranthene		ug/kg ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Benzo(g,h,i)perylene	4700			DNR	<del></del>
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Benzo(k)fluoranthene	4700		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Benzoic Acid	47000		U	DNR	11
	08-21764-NM43KDL	SW8270D	Benzyl Alcohol	47000		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	bis(2-Chloroethoxy) Methane	4700		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Bis-(2-Chloroethyl) Ether	4700		Ü	DNR	- 11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	bis(2-Ethylhexyl)phthalate		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Butylbenzylphthalate	4700		U	DNR	<del>-    </del>
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Carbazole	4700		U	DNR	<del></del>
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Chrysene	12000		H	DNR	<del>- 11</del>
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Dibenz(a,h)anthracene	4700		U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Diethylphthalate	4700		Ü	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Dimethylphthalate	4700		Ü	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Di-n-Butylphthalate	4700		Ü	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Di-n-Octyl phthalate	4700		Ü	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Hexachlorobenzene	4700		Ü	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Hexachlorobutadiene	4700	<u> </u>	Ü	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Hexachlorocyclopentadiene	24000		Ü	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Hexachloroethane	4700		<del>- Ü</del>	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Indeno(1,2,3-cd)pyrene	4700		Ü	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Isophorone	4700		Ü	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Nitrobenzene	4700		Ū	DNR	11

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	DV Qual RC
Q1-D-15.0	08-21764-NM43KDL	SW8270D	N-Nitroso-Di-N-Propylamine		ug/kg	U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	N-Nitrosodiphenylamine		ug/kg	Ü	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Pentachlorophenol		ug/kg	T U	DNR	11
Q1-D-15.0	08-21764-NM43KDL	SW8270D	Phenol		ug/kg	T U	DNR	11
Q1-D-23.0	08-21765-NM43L	NWTPH-Dx	Diesel Range Hydrocarbons		mg/kg	<del>                                     </del>	J	9
Q1-D-23.0	08-21765-NM43L	NWTPH-Dx	Motor Oil		mg/kg	<u> </u>	<del>                                     </del>	9
Q1-D-23.0	08-21765-NM43L	SW8270D	1-Methylnaphthalene	12000	ua/ka	E	DNR	20
Q1-D-23.0	08-21765-NM43L	SW8270D	2-Methylnaphthalene	17000		Ē	DNR	20
Q1-D-23.0	08-21765-NM43L	SW8270D	Acenaphthene	22000		E	DNR	20
Q1-D-23.0	08-21765-NM43L	SW8270D	Anthracene	11000		Ē	DNR	20
Q1-D-23.0	08-21765-NM43L	SW8270D	Benzo(a)anthracene	6500	ug/kg	E	DNR	20
Q1-D-23.0	08-21765-NM43L	SW8270D	Benzo(a)pyrene		ug/kg	<del> </del>	J J	9
Q1-D-23.0	08-21765-NM43L	SW8270D	Benzo(b)fluoranthene	1600	ug/kg	<del>                                     </del>	J	9
Q1-D-23.0	08-21765-NM43L	SW8270D	Benzo(g,h,i)perylene		ug/kg	<del>                                     </del>	J	9
Q1-D-23.0	08-21765-NM43L	SW8270D	Chrysene		ug/kg	E	DNR	20
Q1-D-23.0	08-21765-NM43L	SW8270D	Fluoranthene	48000		E	DNR	20
Q1-D-23.0	08-21765-NM43L	SW8270D	Fluorene	18000		E	DNR	20
Q1-D-23.0	08-21765-NM43L	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg	†	J	9
Q1-D-23.0	08-21765-NM43L	SW8270D	Phenanthrene	120000		ES	DNR	20
Q1-D-23.0	08-21765-NM43L	SW8270D	Pyrene	33000		E	DNR	20
Q1-D-23.0	08-21765-NM43LDL	SW8270D	1,2,4-Trichlorobenzene		ug/kg	T Ū	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	1,2-Dichlorobenzene		ug/kg	ŤŪ	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	1,3-Dichlorobenzene		ug/kg	TŪ	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	1,4-Dichlorobenzene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2,2'-Oxybis(1-Chloropropane)		ug/kg	Ū	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2,4,5-Trichlorophenol		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2,4,6-Trichlorophenol		ug/kg	Ū	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2,4-Dichlorophenol		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2,4-Dimethylphenol		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2,4-Dinitrophenol	20000		U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2,4-Dinitrotoluene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2,6-Dinitrotoluene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2-Chloronaphthalene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2-Chlorophenol		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2-Methylphenol	2000	ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2-Nitroaniline	9900	ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	2-Nitrophenol	2000	ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	3,3'-Dichlorobenzidine	9900	ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	3-Nitroaniline	9900	ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	4,6-Dinitro-2-Methylphenol	20000	ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	4-Bromophenyl-phenylether	2000	ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	4-Chloro-3-methylphenol		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	4-Chloroaniline	9900	ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	4-Chlorophenyl-phenylether	2000	ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	4-Methylphenol	2000	ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	4-Nitroaniline	9900		U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	4-Nitrophenol		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Acenaphthylene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Benzo(a)anthracene		ug/kg		J	9
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Benzo(a)pyrene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Benzo(b)fluoranthene	2100			DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Benzo(g,h,i)perylene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Benzo(k)fluoranthene	2000	ug/kg	U	DNR	11

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	DV Qual RC
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Benzoic Acid	20000		U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Benzyl Alcohol		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	bis(2-Chloroethoxy) Methane		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Bis-(2-Chloroethyl) Ether		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	bis(2-Ethylhexyl)phthalate		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Butylbenzylphthalate		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Carbazole		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Chrysene		ug/kg		J	9
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Dibenz(a,h)anthracene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Dibenzofuran		ug/kg		DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Diethylphthalate		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Dimethylphthalate		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Di-n-Butylphthalate		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Di-n-Octyl phthalate		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Fluoranthene	51000		<u> </u>	J	9
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Hexachlorobenzene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Hexachlorobutadiene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Hexachlorocyclopentadiene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Hexachloroethane		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Isophorone		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Naphthalene		ug/kg	<u> </u>	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Nitrobenzene		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	N-Nitroso-Di-N-Propylamine		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	N-Nitrosodiphenylamine		ug/kg	U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Pentachlorophenol		ug/kg	U U	DNR	11
Q1-D-23.0	08-21765-NM43LDL	SW8270D	Phenoi		ug/kg	U	DNR	11
Q1-D-23,0 SD-2	08-21765-NM43LDL 08-21767-NM43N	SW8270D NWTPH-Dx	Pyrene	52000		<u> </u>	J	9
SD-2 SD-2	08-21767-NM43N	NWTPH-DX	Diesel Range Hydrocarbons Motor Oil		mg/kg	<u> </u>	J J	9
SD-2 SD-2	08-21767-NM43N	SW8270D	1-Methylnaphthalene	11000	mg/kg	<del>                                     </del>	DNR	20
SD-2	08-21767-NM43N	SW8270D	2-Methylnaphthalene	15000		E E	DNR	20
SD-2	08-21767-NM43N	SW8270D	Acenaphthene	17000		E	DNR	20
SD-2	08-21767-NM43N	SW8270D	Anthracene	8800		E	DNR	20
SD-2	08-21767-NM43N	SW8270D	Benzo(a)anthracene		ug/kg ug/kg	<del>                                     </del>	J	9
SD-2	08-21767-NM43N	SW8270D	Benzo(a)pyrene		ug/kg	+	J	9
SD-2	08-21767-NM43N	SW8270D	Benzo(b)fluoranthene		ug/kg ug/kg	<del>  </del>	J	9
SD-2	08-21767-NM43N	SW8270D	Benzo(g,h,i)perylene		ug/kg ug/kg	-	J	9
SD-2	08-21767-NM43N	SW8270D	Chrysene		ug/kg			9
SD-2	08-21767-NM43N	SW8270D	Fluoranthene	26000		E	DNR	20
SD-2	08-21767-NM43N	SW8270D	Fluorene	12000		E	DNR	20
SD-2	08-21767-NM43N	SW8270D	Indeno(1,2,3-cd)pyrene	<u> </u>	ug/kg	_	J	9
SD-2	08-21767-NM43N	SW8270D	Phenanthrene	62000		ES	DNR	20
SD-2	08-21767-NM43N	SW8270D	Pyrene	21000		E	DNR	20
SD-2	08-21767-NM43NDL	SW8270D	1,2,4-Trichlorobenzene	2200		1 U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	1,2-Dichlorobenzene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	1,3-Dichlorobenzene	2200		Ū	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	1,4-Dichlorobenzene	2200		U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2,2'-Oxybis(1-Chloropropane)	2200		Tu	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2,4,5-Trichlorophenol	11000		U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2,4,6-Trichlorophenol	11000		U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2,4-Dichlorophenol	11000		U	DNR	11_
SD-2	08-21767-NM43NDL	SW8270D	2,4-Dimethylphenol	2200	ug/kg	U	DNR	11

						Lab		
Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag		DV Qual RC
SD-2	08-21767-NM43NDL	SW8270D	2,4-Dinitrophenol		ug/kg	U.	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2,4-Dinitrotoluene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2,6-Dinitrotoluene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2-Chloronaphthalene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2-Chlorophenol		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2-Methylphenol		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2-Nitroaniline	11000		U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	2-Nitrophenol		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	3,3'-Dichlorobenzidine	11000		U U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	3-Nitroaniline	11000		U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	4,6-Dinitro-2-Methylphenol	22000		U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	4-Bromophenyl-phenylether		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	4-Chloro-3-methylphenol	11000		U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	4-Chloroaniline	11000		U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	4-Chlorophenyl-phenylether		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	4-Methylphenol		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	4-Nitroaniline	11000		U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	4-Nitrophenol	11000		U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Acenaphthylene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Benzo(a)anthracene		ug/kg		DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Benzo(a)pyrene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Benzo(b)fluoranthene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Benzo(g,h,i)perylene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Benzo(k)fluoranthene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Benzoic Acid	22000	ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Benzyl Alcohol		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	bis(2-Chloroethoxy) Methane		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Bis-(2-Chloroethyl) Ether		ug/kg	Ü	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	bis(2-Ethylhexyl)phthalate		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Butylbenzylphthalate	2200	ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Carbazole	2200	ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Chrysene	4200	ug/kg		DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Dibenz(a,h)anthracene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Dibenzofuran	5300	ug/kg		DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Diethylphthalate		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Dimethylphthalate		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Di-n-Butylphthalate		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Di-n-Octyl phthalate		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Fluoranthene	30000			J	9
SD-2	08-21767-NM43NDL	SW8270D	Hexachlorobenzene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Hexachlorobutadiene		ug/kg	U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Hexachlorocyclopentadiene	11000		T Ū	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Hexachloroethane		ug/kg	T U	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg	1 0	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Isophorone		ug/kg	1 0	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Naphthalene		ug/kg	<del>                                     </del>	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Nitrobenzene		ug/kg	Tu	DNR	11
SD-2	08-21767-NM43NDL	SW8270D	N-Nitroso-Di-N-Propylamine	11000		Ü	DNR	<del>- 11</del>
SD-2	08-21767-NM43NDL	SW8270D	N-Nitrosodiphenylamine		ug/kg		DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Pentachlorophenol	11000			DNR	11
SD-2	08-21767-NM43NDL	SW8270D	Phenol		ug/kg		DNR	11
	08-21767-NM43NDL	SW8270D	Pyrene	31000		+	J	9
SD-2	1130-71/D/*NUM/L3NUM							

						Lab		
Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	DV Qual RC
Q2-A-1.0	08-21714-NM40A	NWTPH-Dx	Diesel Range Hydrocarbons		mg/kg	4	J	8,9
Q2-A-1.0	08-21714-NM40A	SW8270D	Benzoic Acid		ug/kg	U	UJ	10
Q2-A-5.0	08-21715-NM40B	SW8270D	Benzoic Acid		ug/kg	U	UJ	10
Q2-B-1.0	08-21716-NM40C	SW8270D	Benzoic Acid		ug/kg	U	UJ	10
Q2-B-5.0	08-21717-NM40D	SW8270D	Benzoic Acid		ug/kg	U	UJ	10
Q2-C-3.5	08-21718-NM40E	SW8270D	Benzo(a)pyrene		ug/kg	E	DNR	20
Q2-C-3.5	08-21718-NM40E	SW8270D	Benzo(b)fluoranthene		ug/kg	E	DNR	20
Q2-C-3.5	08-21718-NM40E	SW8270D	Benzoic Acid		ug/kg	U	UJ	10
Q2-C-3.5	08-21718-NM40E	SW8270D	Chrysene		ug/kg	E	DNR	20
Q2-C-3.5	08-21718-NM40E	SW8270D	Fluoranthene		ug/kg	E	DNR	20
Q2-C-3.5	08-21718-NM40E	SW8270D	Pyrene		ug/kg	E	DNR	20
Q2-C-3.5	08-21718-NM40EDL	SW8270D	1,2,4-Trichlorobenzene		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	1,2-Dichlorobenzene		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	1,3-Dichlorobenzene		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	1,4-Dichlorobenzene		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	1-Methylnaphthalene		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2,2'-Oxybis(1-Chloropropane)		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2,4,5-Trichlorophenol		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2,4,6-Trichlorophenol		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2,4-Dichlorophenol		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2,4-Dimethylphenol		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2,4-Dinitrophenol		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2,4-Dinitrotoluene		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2,6-Dinitrotoluene		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2-Chloronaphthalene		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2-Chlorophenol		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2-Methylnaphthalene		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2-Methylphenol		ug/kg	Ü	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2-Nitroaniline		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	2-Nitrophenol		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	3,3'-Dichlorobenzidine		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	3-Nitroaniline		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	4,6-Dinitro-2-Methylphenol		ug/kg	U	DNR	11 11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	4-Bromophenyl-phenylether		ug/kg	<del> </del>	DNR	
Q2-C-3.5	08-21718-NM40EDL 08-21718-NM40EDL	SW8270D SW8270D	4-Chloro-3-methylphenol 4-Chloroaniline		ug/kg ug/kg	U	DNR DNR	11 11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	4-Chlorophenyl-phenylether		ug/kg ug/kg	U	DNR	11
Q2-C-3.5 Q2-C-3.5	08-21718-NM40EDL	SW8270D	4-Methylphenol		ug/kg ug/kg	Ü	DNR	11
	08-21718-NM40EDL	SW8270D	4-Nitroaniline		ug/kg ug/kg	l ü	DNR	11
Q2-C-3.5 Q2-C-3.5	08-21718-NM40EDL	SW8270D	4-Nitrophenol		ug/kg ug/kg	U	DNR	11
Q2-C-3.5 Q2-C-3.5	08-21718-NM40EDL	SW8270D	Acenaphthene		ug/kg ug/kg	U	DNR	11
Q2-C-3.5 Q2-C-3.5	08-21718-NM40EDL	SW8270D	Acenaphthylene		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Anthracene		ug/kg	<del>                                     </del>	DNR	11
Q2-C-3.5 Q2-C-3.5	08-21718-NM40EDL	SW8270D	Benzo(a)anthracene		ug/kg ug/kg	<del>  </del>	DNR	11
Q2-C-3.5 Q2-C-3.5	08-21718-NM40EDL	SW8270D	Benzo(g,h,i)perylene		ug/kg	<del>  </del>	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Benzo(k)fluoranthene		ug/kg	<del>  </del>	DNR	11
Q2-C-3.5 Q2-C-3.5	08-21718-NM40EDL	SW8270D	Benzoic Acid		ug/kg	U	DNR	11
Q2-C-3.5 Q2-C-3.5	08-21718-NM40EDL	SW8270D	Benzyl Alcohol		ug/kg ug/kg	l ü	DNR	11
Q2-C-3.5 Q2-C-3.5	08-21718-NM40EDL	SW8270D	bis(2-Chloroethoxy) Methane		ug/kg		DNR	11
Q2-C-3.5 Q2-C-3.5	08-21718-NM40EDL	SW8270D	Bis-(2-Chloroethyl) Ether		ug/kg ug/kg	l ü l	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	bis(2-Ethylhexyl)phthalate		ug/kg ug/kg	$\frac{0}{0}$	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Butylbenzylphthalate		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Carbazole		ug/kg	U	DNR	11
QZ-0-0.0	TOO ETT TO TAIN TOLDE	01102100	σαιραέσιο	010	<del>ugmy</del>		D:11\	

Sample ID	Lab Sample ID	Method	Analyta	Result	Units	Lab Flag	DV Ougl	DV Qual RC
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Analyte Dibenz(a,h)anthracene		ug/kg	Flag	DV Qual	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Dibenzofuran		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Diethylphthalate		ug/kg ug/kg	Ü	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Dimethylphthalate		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Di-n-Butylphthalate		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Di-n-Octyl phthalate		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Fluorene		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Hexachlorobenzene		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Hexachlorobutadiene		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Hexachlorocyclopentadiene		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Hexachloroethane		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg ug/kg		DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Isophorone		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Naphthalene		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Nitrobenzene		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	N-Nitroso-Di-N-Propylamine		ug/kg	Ü	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	N-Nitrosodiphenylamine		ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Pentachlorophenol		ug/kg ug/kg	U	DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Phenanthrene		ug/kg		DNR	11
Q2-C-3.5	08-21718-NM40EDL	SW8270D	Phenol		ug/kg	U	DNR	11
Q2-C-5.0	08-21719-NM40F	SW8270D	Benzoic Acid		ug/kg	Ü	UJ	10
Q2-C-13.0	08-21720-NM40G	SW8270D	1-Methylnaphthalene	51000		E	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	2-Methylnaphthalene	98000		Ē	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Acenaphthene	39000		E	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Anthracene	29000		E	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Benzo(a)anthracene	48000		E	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Benzo(a)pyrene	34000		E	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Benzo(b)fluoranthene	41000		E	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Benzo(k)fluoranthene	25000		E	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Benzoic Acid	2400		<del></del>	UJ	10
Q2-C-13.0	08-21720-NM40G	SW8270D	Carbazole	23000		Ē	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Chrysene	28000		Ē	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Dibenzofuran	44000		Ē	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Fluoranthene	170000		Ē	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Fluorene	68000		E	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Naphthalene	190000		E	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Phenanthrene	230000		Ē	DNR	20
Q2-C-13.0	08-21720-NM40G	SW8270D	Pyrene	100000		Ē	DNR	20
Q2-C-13.0	08-21720-NM40GDL	SW8270D	1,2,4-Trichlorobenzene	4800		- <del>-</del>	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	1,2-Dichlorobenzene	4800		<del>Ŭ</del>	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	1,3-Dichlorobenzene	4800		Ü	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	1,4-Dichlorobenzene	4800		Ū	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2,2'-Oxybis(1-Chloropropane)	4800		Ū	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2,4,5-Trichlorophenol	24000	<u> </u>	Ū	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2,4,6-Trichlorophenol	24000		Ü	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2,4-Dichlorophenol	24000		Ü	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2,4-Dimethylphenol	4800		Ü	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2,4-Dinitrophenol	48000		Ū	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2,4-Dinitrotoluene	24000		Ü	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2,6-Dinitrotoluene	24000		Ü	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2-Chloronaphthalene	4800		Ū	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2-Chlorophenol	4800		Ū	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2-Methylphenol	4800		Ū	DNR	11

Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Lab Flag	DV Oual	DV Qual RC
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2-Nitroaniline		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	2-Nitrophenol		ug/kg	1 0	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	3,3'-Dichlorobenzidine		ug/kg	<del>                                     </del>	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	3-Nitroaniline		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	4,6-Dinitro-2-Methylphenol		ug/kg	<del>                                     </del>	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	4-Bromophenyl-phenylether		ug/kg	<del>                                     </del>	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	4-Chloro-3-methylphenol		ug/kg	<del>  U</del>	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	4-Chloroaniline		ug/kg	<del>                                     </del>	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	4-Chlorophenyl-phenylether		ug/kg	<del>  U</del>	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	4-Methylphenol		ug/kg	1 0	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	4-Nitroaniline		ug/kg	<del>                                     </del>	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	4-Nitrophenol	24000		U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Acenaphthylene		ug/kg ug/kg	1 0	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Benzo(g,h,i)perylene		ug/kg ug/kg	<del>                                     </del>	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Benzoic Acid	48000	<u> </u>	1 11	DNR	11
Q2-C-13.0 Q2-C-13.0	08-21720-NM40GDL	SW8270D	Benzyl Alcohol			U		11
	08-21720-NM40GDL	SW8270D			ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL		bis(2-Chloroethoxy) Methane Bis-(2-Chloroethyl) Ether		ug/kg	U	DNR	
Q2-C-13.0		SW8270D			ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	bis(2-Ethylhexyl)phthalate		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Butylbenzylphthalate		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Dibenz(a,h)anthracene		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Diethylphthalate		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Dimethylphthalate		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Di-n-Butylphthalate		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Di-n-Octyl phthalate		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Hexachlorobenzene		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Hexachlorobutadiene		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Hexachlorocyclopentadiene	24000		U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Hexachloroethane		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Indeno(1,2,3-cd)pyrene	11000		ļ	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Isophorone		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Nitrobenzene		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	N-Nitroso-Di-N-Propylamine	24000		U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	N-Nitrosodiphenylamine		ug/kg	U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Pentachlorophenol	24000		U	DNR	11
Q2-C-13.0	08-21720-NM40GDL	SW8270D	Phenol		ug/kg	U	DNR	11
Q2-C-25.0	08-21721-NM40H	SW8270D	Benzoic Acid		ug/kg	U	UJ	10
Q2-D-3.5	08-21722-NM40I	SW6010B	Copper		mg/kg		J	8
Q2-D-3.5	08-21722-NM40I	SW6010B	Zinc		mg/kg		J	8,9
Q2-D-3.5	08-21722-NM40I	SW8081	delta-BHC		ug/kg	Y	U	22
Q2-D-3.5	08-21722-NM40I	SW8270D	Benzo(a)anthracene	42000		E	DNR	20
Q2-D-3.5	08-21722-NM40I	SW8270D	Benzo(a)pyrene	65000		E	DNR	20
Q2-D-3.5	08-21722-NM40I	SW8270D	Benzo(b)fluoranthene	100000		E	DNR	20
Q2-D-3.5	08-21722-NM40I	SW8270D	Benzo(g,h,i)perylene	34000		E	DNR	20
Q2-D-3.5	08-21722-NM40I	SW8270D	Benzoic Acid		ug/kg	U	UJ	10
Q2-D-3.5	08-21722-NM40I	SW8270D	Chrysene	48000		E	DNR	20
Q2-D-3.5	08-21722-NM40I	SW8270D	Fluoranthene	29000		E	DNR	20
Q2-D-3.5	08-21722-NM40I	SW8270D	Indeno(1,2,3-cd)pyrene	31000		E	DNR	20
Q2-D-3.5	08-21722-NM40l	SW8270D	Pyrene	44000		E	DNR	20
Q2-D-3.5	08-21722-NM40IDL	SW8081	4,4'-DDD	320	ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	4,4'-DDE	320	ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	4,4'-DDT	320	ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Aldrin	160	ug/kg	U	DNR	11

Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Lab Flag	DV Ougl	DV Qual RC
Q2-D-3.5	08-21722-NM40IDL	SW8081	alpha Chlordane		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	alpha-BHC		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	beta-BHC		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	delta-BHC		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Dieldrin		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Endosulfan i		ug/kg ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Endosulfan II		ug/kg	<del>                                     </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Endosulfan Sulfate		ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Endrin		ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Endrin Aldehyde	·	ug/kg	<del>                                     </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Endrin Ketone		ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	gamma Chlordane		ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	gamma-BHC (Lindane)		ug/kg	<del>                                     </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Heptachlor		ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Heptachlor Epoxide		ug/kg	1-5	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Methoxychlor		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8081	Toxaphene	16000	ug/kg ug/kg	<del>                                     </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	1,2,4-Trichlorobenzene		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	1,2-Dichlorobenzene		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	1,3-Dichlorobenzene		ug/kg ug/kg	<del>                                     </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	1,4-Dichlorobenzene		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	1-Methylnaphthalene		ug/kg ug/kg	+	DNR	11
Q2-D-3.5 Q2-D-3.5	08-21722-NM40IDL	SW8270D	2,2'-Oxybis(1-Chloropropane)		ug/kg ug/kg	<del>                                     </del>	DNR	11
Q2-D-3.5 Q2-D-3.5	08-21722-NM40IDL	SW8270D	2,4,5-Trichlorophenol		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5 Q2-D-3.5	08-21722-NM40IDL	SW8270D	2,4,6-Trichlorophenol		ug/kg ug/kg	<del>                                     </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	2,4-Dichlorophenol		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5 Q2-D-3.5	08-21722-NM40IDL	SW8270D	2,4-Dimethylphenol		ug/kg ug/kg	<del>                                     </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	2,4-Dinitrophenol	18000		1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	2,4-Dinitrotoluene		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	2,6-Dinitrotoluene		ug/kg ug/kg	1 0	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	2-Chloronaphthalene		ug/kg	T U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	2-Chlorophenol		ug/kg ug/kg	<del>  U</del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	2-Methylnaphthalene		ug/kg	┼——	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	2-Methylphenol		ug/kg	<del>                                     </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	2-Nitroaniline		ug/kg	Ü	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	2-Nitrophenol		ug/kg	Ū	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	3,3'-Dichlorobenzidine		ug/kg	<del>l ű</del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	3-Nitroaniline		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	4,6-Dinitro-2-Methylphenol	18000		Ü	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	4-Bromophenyl-phenylether		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	4-Chloro-3-methylphenol		ug/kg	Ü	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	4-Chloroaniline		ug/kg	Ü	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	4-Chlorophenyl-phenylether		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	4-Methylphenol		ug/kg	Ü	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	4-Nitroaniline		ug/kg	Ü	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	4-Nitrophenol		ug/kg	T U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Acenaphthene		ug/kg	<del>                                     </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Acenaphthylene		ug/kg	<del> </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Anthracene	10000			DNR	<del>- ;;</del> -
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Benzo(k)fluoranthene	47000		<del>                                     </del>	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Benzoic Acid	18000		U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Benzyl Alcohol		ug/kg	Ū	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	bis(2-Chloroethoxy) Methane		ug/kg	T Ü	DNR	11

Sample ID	1			1		Lab		İ
	Lab Sample ID	Method	Analyte	Result	Units	Flag		DV Qual RC
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Bis-(2-Chloroethyl) Ether		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	bis(2-Ethylhexyl)phthalate		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Butylbenzylphthalate		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Carbazole		ug/kg		DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Dibenz(a,h)anthracene		ug/kg		DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Dibenzofuran		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Diethylphthalate		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Dimethylphthalate		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Di-n-Butyiphthalate		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Di-n-Octyl phthalate		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Fluorene		ug/kg		DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Hexachlorobenzene	1800	ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Hexachlorobutadiene	1800	ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Hexachlorocyclopentadiene	9100	ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Hexachloroethane	1800	ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Isophorone	1800	ug/kg	U	DNR	11
	08-21722-NM40IDL	SW8270D	Naphthalene		ug/kg		DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Nitrobenzene		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	N-Nitroso-Di-N-Propylamine	9100	ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	N-Nitrosodiphenylamine		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Pentachlorophenol		ug/kg	U	DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Phenanthrene	24000			DNR	11
Q2-D-3.5	08-21722-NM40IDL	SW8270D	Phenol		ug/kg	U	DNR	11
	08-21723-NM40J	SW6010B	Copper		mg/kg		J	8
	08-21723-NM40J	SW6010B	Zinc		mg/kg		J	8,9
	08-21723-NM40J	SW8081	Aldrin		ug/kg	Y	U	22
	08-21723-NM40J	SW8081	delta-BHC		ug/kg	Y	U	22
	08-21723-NM40J	SW8081	gamma Chlordane		ug/kg	Y	U	22
	08-21723-NM40J	SW8082	Aroclor 1248		ug/kg	Y	U	22
	08-21723-NM40J	SW8270D	Benzo(a)anthracene	250000	ua/ka	E	DNR	20
	08-21723-NM40J	SW8270D	Benzo(a)pyrene	190000		E	DNR	20
	08-21723-NM40J	SW8270D	Benzo(b)fluoranthene	280000		E	DNR	20
	08-21723-NM40J	SW8270D	Benzo(g,h,i)perylene	76000		E	DNR	20
	08-21723-NM40J	SW8270D	Benzo(k)fluoranthene	150000		<del>                                     </del>	DNR	20
	08-21723-NM40J	SW8270D	Benzoic Acid		ug/kg	1 0	UJ	10
	08-21723-NM40J	SW8270D	Chrysene	220000		T E	DNR	20
	08-21723-NM40J	SW8270D	Fluoranthene	170000		TE	DNR	20
	08-21723-NM40J	SW8270D	Indeno(1,2,3-cd)pyrene	83000		E	DNR	20
	08-21723-NM40J	SW8270D	Pyrene	270000		E	DNR	20
	08-21723-NM40JDL	SW8081	4,4'-DDD		ug/kg	1 0	DNR	11
	08-21723-NM40JDL	SW8081	4,4'-DDE		ug/kg	1 0	DNR	11
	08-21723-NM40JDL	SW8081	4,4'-DDT		ug/kg	1 0 1	DNR	11
	08-21723-NM40JDL	SW8081	Aldrin		ug/kg	1 0	DNR	11
	08-21723-NM40JDL	SW8081	alpha Chlordane		ug/kg	<del>l ŭ l</del>	DNR	11
	08-21723-NM40JDL	SW8081	alpha-BHC		ug/kg ug/kg	1 0 1	DNR	11
	08-21723-NM40JDL	SW8081	beta-BHC		ug/kg		DNR	<del></del>
	08-21723-NM40JDL	SW8081	delta-BHC		ug/kg	1 0 1	DNR	11
	08-21723-NM40JDL	SW8081	Dieldrin		ug/kg	U	DNR	11
	08-21723-NM40JDL	SW8081	Endosulfan I		ug/kg ug/kg	1 0 1	DNR	11
	08-21723-NM40JDL	SW8081	Endosulfan II		ug/kg ug/kg		DNR	11
	08-21723-NM40JDL	SW8081	Endosulfan Sulfate		ug/kg ug/kg	1 0 1	DNR	11
	08-21723-NM40JDL	SW8081	Endrin		ug/kg ug/kg	1 0 1	DNR	11
<b>ਘ∠⁻レ⁻∪.∪  </b>	08-21723-NM40JDL	SW8081	Endrin Aldehyde		ug/kg ug/kg	1 0	DNR	11

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag		DV Qual RC
Q2-D-5.0	08-21723-NM40JDL	SW8081	Endrin Ketone		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8081	gamma Chlordane		ug/kg_	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8081	gamma-BHC (Lindane)		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8081	Heptachlor		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8081	Heptachlor Epoxide		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8081	Methoxychlor		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8081	Toxaphene	16000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	1,2,4-Trichlorobenzene		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	1,2-Dichlorobenzene		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	1,3-Dichlorobenzene		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	1,4-Dichlorobenzene		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	1-Methylnaphthalene		ug/kg	U_	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2,2'-Oxybis(1-Chloropropane)		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2,4,5-Trichlorophenol	38000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2,4,6-Trichlorophenol	38000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2,4-Dichlorophenol	38000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2,4-Dimethylphenol		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2,4-Dinitrophenol	76000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2,4-Dinitrotoluene	38000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2,6-Dinitrotoluene	38000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2-Chloronaphthalene	7600	ug/kg	C	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2-Chlorophenol	7600	ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2-Methylnaphthalene	7600	ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2-Methylphenol	7600	ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2-Nitroaniline	38000	ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	2-Nitrophenol	7600	ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	3,3'-Dichlorobenzidine	38000	ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	3-Nitroaniline	38000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	4,6-Dinitro-2-Methylphenol	76000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	4-Bromophenyl-phenylether		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	4-Chloro-3-methylphenol	38000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	4-Chloroaniline	38000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	4-Chlorophenyl-phenylether		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	4-Methylphenol		ug/kg	U	DNR	
Q2-D-5.0	08-21723-NM40JDL	SW8270D	4-Nitroaniline	38000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	4-Nitrophenol	38000			DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Acenaphthene	10000		1	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Acenaphthylene	10000			DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Anthracene	22000			DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Benzoic Acid	76000		U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Benzyl Alcohol		ug/kg	<del>                                     </del>	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	bis(2-Chloroethoxy) Methane		ug/kg	1 0	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Bis-(2-Chloroethyl) Ether	7600		<del>                                     </del>	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	bis(2-Ethylhexyl)phthalate		ug/kg	T Ü	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Butylbenzylphthalate		ug/kg	1 0	DNR	<del></del>
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Carbazole	7600		<del>                                     </del>	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Dibenz(a,h)anthracene	29000		┯┷	DNR	<del></del>
Q2-D-5.0 Q2-D-5.0	08-21723-NM40JDL	SW8270D	Dibenzofuran	7600		1-0-1	DNR	<del></del>
Q2-D-5.0 Q2-D-5.0	08-21723-NM40JDL	SW8270D	Diethylphthalate	7600		1 0	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Dimethylphthalate	7600		1 0	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Di-n-Butylphthalate	7600		1 0	DNR	<del></del>
Q2-D-5.0 Q2-D-5.0	08-21723-NM40JDL	SW8270D	Di-n-Octyl phthalate	7600			DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Fluorene	7600		1 0	DNR	11

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	<del></del>	DV Qual RC
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Hexachlorobenzene		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Hexachlorobutadiene		ug/kg	- U	DNR	11
Q2-D-5.0	08-21723-NM40JDL 08-21723-NM40JDL	SW8270D SW8270D	Hexachlorocyclopentadiene	38000		U	DNR	11
Q2-D-5.0			Hexachloroethane		ug/kg	U	DNR	11
Q2-D-5.0	08-21723-NM40JDL	SW8270D	Isophorone		ug/kg	U	DNR	11
Q2-D-5.0 Q2-D-5.0	08-21723-NM40JDL	SW8270D	Naphthalene		ug/kg	<u> </u>	DNR	11
Q2-D-5.0 Q2-D-5.0	08-21723-NM40JDL	SW8270D	Nitrobenzene		ug/kg	U	DNR	11
Q2-D-5.0 Q2-D-5.0	08-21723-NM40JDL 08-21723-NM40JDL	SW8270D SW8270D	N-Nitroso-Di-N-Propylamine	38000		U	DNR	11
Q2-D-5.0 Q2-D-5.0	08-21723-NM40JDL	SW8270D	N-Nitrosodiphenylamine		ug/kg	U	DNR	11
	08-21723-NM40JDL	SW8270D	Pentachlorophenol	38000		U	DNR	11
Q2-D-5.0 Q2-D-5.0	08-21723-NM40JDL	SW8270D	Phenanthrene		ug/kg	U	DNR	11 11
Q2-D-5.0 Q2-D-10.0	08-21723-NM40JDL 08-21724-NM40K	SW8270D SW8270D	Phenol		ug/kg	U	DNR	20
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	1-Methylnaphthalene	25000		E	DNR	20
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	2-Methylnaphthalene	54000		E	DNR	20
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	Acenaphthene	13000		E	DNR	
	08-21724-NM40K	SW8270D	Anthracene	11000		E	DNR	20 20
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	Benzo(a)anthracene		ug/kg	E	DNR	20
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	Benzo(b)fluoranthene		ug/kg	E	DNR	
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	Benzoic Acid		ug/kg	U	UJ	10 20
	08-21724-NM40K	SW8270D	Carbazole		ug/kg	E	DNR	
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	Chrysene Dibenzofuran		ug/kg	E	DNR	20 20
	08-21724-NM40K	SW8270D		15000			DNR	
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	Fluoranthene Fluorene	63000		I E	DNR	20 20
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	<del></del>	25000		E E	DNR	20
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	Naphthalene Phenanthrene	54000 120000		ES	DNR DNR	20
Q2-D-10.0 Q2-D-10.0	08-21724-NM40K	SW8270D	Pyrene	25000		E	DNR	20
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	1,2,4-Trichlorobenzene		ug/kg ug/kg	<u> </u>	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	1,2-Dichlorobenzene		ug/kg ug/kg	1 0	DNR	<u>'                                 </u>
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	1,3-Dichlorobenzene		ug/kg ug/kg	1 0	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	1,4-Dichlorobenzene	2400	ug/kg ug/kg	1 0	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	2,2'-Oxybis(1-Chloropropane)		ug/kg ug/kg	<del>  0</del>	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	2,4,5-Trichlorophenol	12000		1 0	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	2,4,6-Trichlorophenol	12000		1 5	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	2,4-Dichlorophenol	12000		1 0	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	2,4-Dimethylphenol	2400		1 0	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	2,4-Dinitrophenol	24000			DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	2,4-Dinitrotoluene	12000			DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	2,6-Dinitrotoluene	12000			DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	2-Chloronaphthalene	2400		Ü	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	2-Chlorophenol	2400		T U	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	2-Methylphenol	2400		1 0	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	2-Nitroaniline	12000		<del>                                     </del>	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	2-Nitrophenol	2400			DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	3,3'-Dichlorobenzidine	12000		1 0	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	3-Nitroaniline	12000		1 0	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	4,6-Dinitro-2-Methylphenol	24000		T U	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	4-Bromophenyl-phenylether	2400		T Ü	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	4-Chloro-3-methylphenol	12000		T Ü	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	4-Chloroaniline	12000		T Ü	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	4-Chlorophenyl-phenylether	2400		<del>                                     </del>	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	4-Methylphenol	2400		<del>                                     </del>	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	4-Nitroaniline	12000		<del>1 ਹੈ  </del>	DNR	11

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag		DV Qual RC
Q2-D-10.0	08-21724-NM40KDL	SW8270D	4-Nitrophenol		ug/kg	U	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	Acenaphthylene		ug/kg	U	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	Benzo(a)pyrene		ug/kg		DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	Benzo(g,h,i)perylene		ug/kg	U	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	Benzo(k)fluoranthene		ug/kg		DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	Benzoic Acid	24000		U	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	Benzyl Alcohol		ug/kg	U	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	bis(2-Chloroethoxy) Methane		ug/kg	U	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D SW8270D	Bis-(2-Chloroethyl) Ether		ug/kg	U	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL		bis(2-Ethylhexyl)phthalate		ug/kg	U	DNR	11
	08-21724-NM40KDL	SW8270D	Butylbenzylphthalate		ug/kg	U	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	Dibenz(a,h)anthracene		ug/kg	U	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D SW8270D	Diethylphthalate		ug/kg	U	DNR	11 11
	08-21724-NM40KDL 08-21724-NM40KDL		Dimethylphthalate		ug/kg	U	DNR	
Q2-D-10.0	08-21724-NM40KDL	SW8270D	Di-n-Butylphthalate		ug/kg	U	DNR	11 11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D SW8270D	Di-n-Octyl phthalate Hexachlorobenzene		ug/kg ug/kg	U	DNR DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	Hexachlorobutadiene		ug/kg ug/kg	1 0	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	Hexachlorocyclopentadiene	12000		<del>                                     </del>	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	Hexachloroethane		ug/kg ug/kg	1 0	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg ug/kg	1 0	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	Isophorone		ug/kg ug/kg	1 0	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	Nitrobenzene		ug/kg ug/kg	1 0	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	N-Nitroso-Di-N-Propylamine	12000		1 0	DNR	11
Q2-D-10.0 Q2-D-10.0	08-21724-NM40KDL	SW8270D	N-Nitrosodiphenylamine		ug/kg ug/kg	1 0	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	Pentachlorophenol	12000		<del>  U</del>	DNR	11
Q2-D-10.0	08-21724-NM40KDL	SW8270D	Phenol		ug/kg	1 0	DNR	11
Q2-D-13.0	08-21725-NM40L	SW8270D	2-Methylnaphthalene		ug/kg	E	DNR	20
Q2-D-13.0	08-21725-NM40L	SW8270D	Benzoic Acid		ug/kg	Ū	UJ	10
Q2-D-13.0	08-21725-NM40L	SW8270D	Naphthalene	14000		Ē	DNR	20
Q2-D-13.0	08-21725-NM40L	SW8270D	Phenanthrene		ug/kg	Ē	DNR	20
Q2-D-13.0	08-21725-NM40LDL	SW8270D	1,2,4-Trichlorobenzene		ug/kg	T U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	1,2-Dichlorobenzene		ug/kg	Ū	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	1,3-Dichlorobenzene		ug/kg	Ū	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	1,4-Dichlorobenzene		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	1-Methylnaphthalene		ug/kg		DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2,2'-Oxybis(1-Chloropropane)		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2,4,5-Trichlorophenol	1600	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2,4,6-Trichlorophenol		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2,4-Dichlorophenol	1600	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2,4-Dimethylphenol	320	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2,4-Dinitrophenol	3200	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2,4-Dinitrotoluene	1600	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2,6-Dinitrotoluene	1600	ug/kg	$\Box$	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2-Chloronaphthalene		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2-Chlorophenol		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2-Methylphenol		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2-Nitroaniline		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	2-Nitrophenol		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	3,3'-Dichlorobenzidine	1600		U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	3-Nitroaniline		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	4,6-Dinitro-2-Methylphenol		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	4-Bromophenyl-phenylether	320	ug/kg	U	DNR	11

Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Lab Flag	DV Ousi	DV Qual RC
Q2-D-13.0	08-21725-NM40LDL	SW8270D	4-Chloro-3-methylphenol		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	4-Chloroaniline		ug/kg ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	4-Chlorophenyl-phenylether		ug/kg ug/kg	l ü	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	4-Methylphenol		ug/kg ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	4-Nitroaniline		ug/kg	T U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	4-Nitrophenol		ug/kg	<del>                                     </del>	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Acenaphthene		ug/kg	<del>                                     </del>	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Acenaphthylene		ug/kg	l u	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Anthracene		ug/kg	<del>                                     </del>	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Benzo(a)anthracene		ug/kg	<b>-</b>	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Benzo(a)pyrene		ug/kg	<del> </del>	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Benzo(b)fluoranthene		ug/kg		DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Benzo(g,h,i)perylene		ug/kg	<del> </del>	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Benzo(k)fluoranthene		ug/kg	<b></b>	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Benzoic Acid		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Benzyl Alcohol		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	bis(2-Chloroethoxy) Methane		ug/kg	Ü	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Bis-(2-Chloroethyl) Ether		ug/kg	Ū	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	bis(2-Ethylhexyl)phthalate		ug/kg	Ū	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Butylbenzylphthalate		ug/kg	<del>- i</del>	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Carbazole		ug/kg		DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Chrysene		ug/kg		DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Dibenz(a,h)anthracene		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Dibenzofuran		ug/kg		DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Diethylphthalate		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Dimethylphthalate		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Di-n-Butylphthalate		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Di-n-Octyl phthalate		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Fluoranthene		ug/kg		DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Fluorene		ug/kg		DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Hexachlorobenzene		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Hexachlorobutadiene	320	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Hexachlorocyclopentadiene	1600	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Hexachloroethane	320	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg		DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Isophorone	320	ug/kg	U	DNR	11_
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Nitrobenzene	320	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	N-Nitroso-Di-N-Propylamine		ug/kg	Ü	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	N-Nitrosodiphenylamine	320	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Pentachlorophenol	1600	ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Phenol		ug/kg	U	DNR	11
Q2-D-13.0	08-21725-NM40LDL	SW8270D	Pyrene	2600			DNR	11
Q2-D-18.0	08-21726-NM40M	SW8270D	1-Methylnaphthalene	86000		E	DNR	20
Q2-D-18.0	08-21726-NM40M	SW8270D	2-Methylnaphthalene	180000		E	DNR	20
Q2-D-18.0	08-21726-NM40M	SW8270D	Acenaphthene	39000		E	DNR	20
Q2-D-18.0	08-21726-NM40M	SW8270D	Anthracene	31000		E	DNR	20
Q2-D-18.0	08-21726-NM40M	SW8270D	Benzo(a)anthracene	37000		Е	DNR	20
Q2-D-18.0	08-21726-NM40M	SW8270D	Benzo(a)pyrene	22000		E	DNR	20
Q2-D-18.0	08-21726-NM40M	SW8270D	Benzo(b)fluoranthene	24000		E	DNR	20
Q2-D-18.0	08-21726-NM40M	SW8270D	Benzoic Acid	2100		U	UJ	10
Q2-D-18.0	08-21726-NM40M	SW8270D	Chrysene	22000		E	DNR	20
Q2-D-18.0	08-21726-NM40M	SW8270D	Dibenzofuran	40000		E	DNR	20
Q2-D-18.0	08-21726-NM40M	SW8270D_	Fluoranthene	160000	ug/kg	E	DNR	20

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	<del></del>	DV Qual RC
Q2-D-18.0	08-21726-NM40M	SW8270D	Fluorene		ug/kg	E	DNR	20
Q2-D-18.0	08-21726-NM40M	SW8270D	Naphthalene	200000		E	DNR	20
Q2-D-18.0	08-21726-NM40M 08-21726-NM40M	SW8270D SW8270D	Phenanthrene	290000		ES	DNR	20
Q2-D-18.0	08-21726-NM40MDL		Pyrene 4.2.4 Trichlershanzana	86000		E	DNR	20
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D SW8270D	1,2,4-Trichlorobenzene		ug/kg	U	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D SW8270D	1,2-Dichlorobenzene		ug/kg	U	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D SW8270D	1,3-Dichlorobenzene		ug/kg	U	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D	1,4-Dichlorobenzene		ug/kg	U	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D	2,2'-Oxybis(1-Chloropropane)		ug/kg		DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D	2,4,5-Trichlorophenol	42000		U	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D	2,4,6-Trichlorophenol 2,4-Dichlorophenol	42000		U	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D		42000		U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	2,4-Dimethylphenol 2,4-Dinitrophenol	85000	ug/kg	U	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D				1 0	DNR	11
Q2-D-18.0 Q2-D-18.0		SW8270D SW8270D	2,4-Dinitrotoluene	42000			DNR	
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D SW8270D	2,6-Dinitrotoluene	42000		U	DNR	11
	08-21726-NM40MDL 08-21726-NM40MDL		2-Chloronaphthalene		ug/kg	U	DNR	11 11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D SW8270D	2-Chlorophenol		ug/kg	<del>  U</del>	DNR	
		SW8270D SW8270D	2-Methylphenol 2-Nitroaniline		ug/kg		DNR	11
Q2-D-18.0	08-21726-NM40MDL 08-21726-NM40MDL			42000		U	DNR	11
Q2-D-18.0		SW8270D	2-Nitrophenol		ug/kg	U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	3,3'-Dichlorobenzidine 3-Nitroaniline	42000		U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D		42000		U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	4,6-Dinitro-2-Methylphenol	85000		U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	4-Bromophenyl-phenylether		ug/kg	U	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL 08-21726-NM40MDL	SW8270D SW8270D	4-Chloro-3-methylphenol 4-Chloroaniline	42000 42000		U	DNR DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D			ug/kg ug/kg	1 0	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D	4-Chlorophenyl-phenylether		ug/kg ug/kg	1 0	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D	4-Methylphenol 4-Nitroaniline	42000		1 0	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D	4-Nitrophenol	42000		1 0	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Acenaphthylene		ug/kg ug/kg	1 0	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Benzo(g,h,i)perylene		ug/kg	U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Benzo(k)fluoranthene	21000		<del>                                     </del>	DNR	11
Q2-D-18.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D	Benzoic Acid	85000		T U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Benzyl Alcohol		ug/kg ug/kg		DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	bis(2-Chloroethoxy) Methane		ug/kg ug/kg	1 0	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Bis-(2-Chloroethyl) Ether		ug/kg	U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	bis(2-Ethylhexyl)phthalate		ug/kg ug/kg		DNR	11
Q2-D-10.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D	Butylbenzylphthalate		ug/kg	1-0-	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Carbazole	10000		<del>                                     </del>	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Dibenz(a,h)anthracene		ug/kg ug/kg	U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Diethylphthalate		ug/kg	U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Dimethylphthalate		ug/kg ug/kg	U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Di-n-Butylphthalate	8500		U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Di-n-Octyl phthalate		ug/kg ug/kg	1 0	DNR	11
Q2-D-10.0 Q2-D-18.0	08-21726-NM40MDL	SW8270D	Hexachlorobenzene	8500		U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Hexachlorobutadiene		ug/kg ug/kg	T U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Hexachlorocyclopentadiene	42000			DNR	11
	08-21726-NM40MDL	SW8270D	Hexachloroethane		ug/kg ug/kg		DNR	11
	08-21726-NM40MDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg ug/kg		DNR	11
	08-21726-NM40MDL	SW8270D	Isophorone		ug/kg	U	DNR	11
	08-21726-NM40MDL	SW8270D	Nitrobenzene	8500		U	DNR	11

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	DV Qual RC
Q2-D-18.0	08-21726-NM40MDL	SW8270D	N-Nitroso-Di-N-Propylamine	42000	ug/kg	U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	N-Nitrosodiphenylamine	8500	ug/kg	U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Pentachlorophenol	42000	ug/kg	U	DNR	11
Q2-D-18.0	08-21726-NM40MDL	SW8270D	Phenol	8500	ug/kg	U	DNR	11
Q2-D-35.0	08-21727-NM40N	SW8270D	Benzoic Acid	630	ug/kg	Ü	UJ	10

#### **Quendall Terminal -Transect Q3**

Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Lab Flag	DV Qual	DV Qual RC
Q3-A-1.0	08-21748-NM42A	SW8270D	Fluoranthene	1400000		E	DNR	20
Q3-A-1.0	08-21748-NM42ADL	SW8270D	1,2,4-Trichlorobenzene		ug/kg	1 -	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	1,2-Dichlorobenzene	28000		1 0	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	1,3-Dichlorobenzene	28000		<del>  U</del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	1,4-Dichlorobenzene	28000		1 0	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	1-Methylnaphthalene	28000		1 0	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2,2'-Oxybis(1-Chloropropane)	28000		<del>                                     </del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2,4,5-Trichlorophenol	140000		<del>                                     </del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2,4,6-Trichlorophenol	140000		ΙÜ	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2,4-Dichlorophenol	140000		<del>l ü</del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2,4-Dimethylphenol	28000		<del>l ü</del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2,4-Dinitrophenol	280000		Ι <del>ΰ</del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2,4-Dinitrotoluene	140000		<del>                                     </del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2,6-Dinitrotoluene	140000		1 0	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2-Chloronaphthalene	28000		<del>  Ŭ</del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2-Chlorophenol	28000		<del>  U</del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2-Methylnaphthalene	28000	ug/kg ug/kg	1 0	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2-Methylphenol	28000		1 0	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2-Nitroaniline	140000		T U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	2-Nitrophenol	28000		1 0	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	3,3'-Dichlorobenzidine	140000		T U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	3-Nitroaniline	140000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	4,6-Dinitro-2-Methylphenol	280000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	4-Bromophenyl-phenylether	28000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	4-Chloro-3-methylphenol	140000		Ü	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	4-Chloroaniline	140000		T U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	4-Chlorophenyl-phenylether	28000		Ü	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	4-Methylphenol	28000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	4-Nitroaniline	140000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	4-Nitrophenol	140000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Acenaphthene	88000		+	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Acenaphthylene	28000	ua/ka		DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Anthracene	92000	ua/ka	<del>                                     </del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Benzo(a)anthracene	380000			DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Benzo(a)pyrene	260000			DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Benzo(b)fluoranthene	270000		ļ	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Benzo(g,h,i)perylene	120000			DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Benzo(k)fluoranthene	210000			DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Benzoic Acid	280000	_ <del>```</del>	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Benzyl Alcohol	28000		Ū	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	bis(2-Chloroethoxy) Methane	28000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Bis-(2-Chloroethyl) Ether	28000		Ū	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	bis(2-Ethylhexyl)phthalate	28000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Butylbenzylphthalate	28000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Carbazole	28000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Chrysene	460000			DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Dibenz(a,h)anthracene	28000	<u> </u>	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Dibenzofuran	28000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Diethylphthalate	28000		Ū	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Dimethylphthalate	28000		<del>                                     </del>	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Di-n-Butylphthalate	28000		Ū	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Di-n-Octyl phthalate	28000		Ū	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Fluorene	46000			DNR	11
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### **Quendall Terminal -Transect Q3**

						Lab		
Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	DV Qual RC
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Hexachlorobenzene	28000	ug/kg	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Hexachlorobutadiene	28000	ug/kg	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Hexachlorocyclopentadiene	140000	ug/kg	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Hexachloroethane	28000	ug/kg	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Indeno(1,2,3-cd)pyrene	110000	ug/kg		DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Isophorone	28000	ug/kg	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Naphthalene	28000	ug/kg	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Nitrobenzene	28000	ug/kg	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	N-Nitroso-Di-N-Propylamine	140000	ug/kg	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	N-Nitrosodiphenylamine	28000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Pentachlorophenol	140000	ug/kg	U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Phenanthrene	260000	ug/kg		DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Phenol	28000		U	DNR	11
Q3-A-1.0	08-21748-NM42ADL	SW8270D	Pyrene	970000	ug/kg		DNR	11

#### **Baxter Mills - Transect B1**

Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Lab Flag	DV Qual	DV Qual RC
B1-A-1.0	08-21932-NM59A	NWTPH-Dx	Diesel Range Hydrocarbons		mg/kg		J	9
B1-A-3.5	08-21933-NM59B	SW8270D	Benzyl Alcohol		ug/kg	U	UJ	5B
B1-A-5.0	08-21934-NM59C	SW8270D	Benzyl Alcohol		ug/kg	U	UJ	5B
B1-A-10.0	08-21935-NM59D	SW8270D	Benzyl Alcohol		ug/kg	U	UJ	5B
B1-A-15.0	08-21936-NM59E	SW8270D	Benzyl Alcohol		ug/kg	U	UJ	5B
B1-A-20.0	08-21937-NM59F	SW8270D	Benzyl Alcohol		ug/kg	U	UJ	5B
B1-A-24.0	08-21938-NM59G	SW8270D	Benzyl Alcohol		ug/kg	U	UJ	5B
B1-B-1.0	08-21939-NM59H	SW8270D	Benzo(a)pyrene	55000		E	DNR	20
B1-B-1.0	08-21939-NM59H	SW8270D	Chrysene	51000		E	DNR	20
B1-B-1.0	08-21939-NM59H	SW8270D	Fluoranthene	67000		E	DNR	20
B1-B-1.0	08-21939-NM59H	SW8270D	Phenanthrene	82000		E	DNR	20
B1-B-1.0	08-21939-NM59H	SW8270D	Pyrene	100000		E	DNR	20
B1-B-1.0	08-21939-NM59HDL	SW8270D	1,2,4-Trichlorobenzene		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	1,2-Dichlorobenzene		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	1,3-Dichlorobenzene		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	1,4-Dichlorobenzene		ug/kg	l U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	1-Methylnaphthalene		ug/kg		DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2,2'-Oxybis(1-Chloropropane)		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2,4,5-Trichlorophenol	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2,4,6-Trichlorophenol	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2,4-Dichlorophenol	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2,4-Dimethylphenol		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2,4-Dinitrophenol	24000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2,4-Dinitrotoluene	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2,6-Dinitrotoluene	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2-Chloronaphthalene		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2-Chlorophenol		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2-Methylnaphthalene	10000			DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2-Methylphenol		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2-Nitroaniline	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	2-Nitrophenol		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	3,3'-Dichlorobenzidine	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	3-Nitroaniline	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	4,6-Dinitro-2-Methylphenol	24000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	4-Bromophenyl-phenylether		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	4-Chloro-3-methylphenol	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	4-Chloroaniline	12000	<u> </u>	U	DNR	11
B1-B-1.0	08-21939-NM59HDL		4-Chlorophenyl-phenylether	2400		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	4-Methylphenol	2400		U	DNR	11
B1-B-1.0	08-21939-NM59HDL		4-Nitroaniline	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	4-Nitrophenol	12000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL		Acenaphthene	15000			DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Acenaphthylene	2400		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Anthracene	22000			DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Benzo(a)anthracene	47000			DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Benzo(b)fluoranthene	40000			DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Benzo(g,h,i)perylene	19000			DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Benzo(k)fluoranthene	38000		<del></del>	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Benzoic Acid	24000		U	DNR	11
B1-B-1.0	08-21939-NM59HDL		Benzyl Alcohol	2400		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	bis(2-Chloroethoxy) Methane	2400		U	DNR	11
B1-B-1.0	08-21939-NM59HDL		Bis-(2-Chloroethyl) Ether	2400		U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	bis(2-Ethylhexyl)phthalate	2400	ug/kg	U	DNR	11

### Baxter Mills - Transect B1

	T T	T				Lab	I .	
Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	DV Qual RC
B1-B-1.0	08-21939-NM59HDL	SW8270D	Butylbenzylphthalate	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Carbazole	5600	ug/kg		DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Dibenz(a,h)anthracene	3600	ug/kg		DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Dibenzofuran	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Diethylphthalate	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Dimethylphthalate	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Di-n-Butylphthalate	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Di-n-Octyl phthalate	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Fluorene	8700	ug/kg		DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Hexachlorobenzene	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Hexachlorobutadiene	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Hexachlorocyclopentadiene	12000	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Hexachloroethane		ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Indeno(1,2,3-cd)pyrene	14000	ug/kg		DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Isophorone	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Naphthalene	16000	ug/kg		DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Nitrobenzene	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	N-Nitroso-Di-N-Propylamine	12000	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	N-Nitrosodiphenylamine	2400	ug/kg	U	DNR	11
B1-B-1.0	08-21939-NM59HDL	SW8270D	Pentachlorophenol	12000	ug/kg	U	DNR	11_
B1-B-1.0	08-21939-NM59HDL	SW8270D	Phenol	2400	ug/kg	U	DNR	11
B1-B-3.0	08-21940-NM59I	SW8270D	Benzyl Alcohol	61	ug/kg	U	UJ	5B
DS-1	08-21943-NM59L	NWTPH-Dx	Diesel Range Hydrocarbons	320	mg/kg		J	9

			Quendan Terminar - Bor	1		Lab	T	DV Qual
Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	RC
Q4-2.5	08-29725-NX66A	SW8270D	Anthracene		ug/kg	E	DNR	20
Q4-2.5	08-29725-NX66A	SW8270D	Benzo(a)anthracene		ug/kg	E	DNR	20
Q4-2.5	08-29725-NX66A	SW8270D	Benzo(a)pyrene		ug/kg	T E	DNR	20
Q4-2.5	08-29725-NX66A	SW8270D	Benzo(b)fluoranthene	18000	ug/kg	E	DNR	20
Q4-2.5	08-29725-NX66A	SW8270D	Benzo(k)fluoranthene		ug/kg	TĒ.	DNR	20
Q4-2.5	08-29725-NX66A	SW8270D	Chrysene	67000	ug/kg	E	DNR	20
Q4-2.5	08-29725-NX66A	SW8270D	Fluoranthene	23000	ua/ka	E	DNR	20
Q4-2.5	08-29725-NX66A	SW8270D	Fluorene		ug/kg	E	DNR	20
Q4-2.5	08-29725-NX66A	SW8270D	Naphthalene		ug/kg	E	DNR	20
Q4-2.5	08-29725-NX66A	SW8270D	Phenanthrene		ug/kg	E	DNR	20
Q4-2.5	08-29725-NX66A	SW8270D	Pyrene	32000		E	DNR	20
Q4-2.5	08-29725-NX66ADL	SW8270D	1-Methylnaphthalene		ug/kg		DNR	11
Q4-2.5	08-29725-NX66ADL	SW8270D	2-Methylnaphthalene		ug/kg	<del>                                     </del>	DNR	11
Q4-2.5	08-29725-NX66ADL	SW8270D	Acenaphthene		ug/kg	_	DNR	11
Q4-2.5	08-29725-NX66ADL	SW8270D	Acenaphthylene		ug/kg	U	DNR	11
Q4-2.5	08-29725-NX66ADL	SW8270D	Benzo(g,h,i)perylene		ug/kg		DNR	11
Q4-2.5	08-29725-NX66ADL	SW8270D	Dibenz(a,h)anthracene		ug/kg		DNR	11
Q4-2.5	08-29725-NX66ADL	SW8270D	Dibenzofuran		ug/kg		DNR	11
Q4-2.5	08-29725-NX66ADL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg		DNR	11
Q4-15.0	08-29727-NX66C	SW8021B	Benzene		ug/kg		J	9
Q4-15.0	08-29727-NX66C	SW8270D	Dibenzofuran	130000	ua/ka		J	9
Q4-15.0	08-29727-NX66C	SW8270D	Naphthalene	1100000		<del> </del> E	DNR	20
Q4-15.0	08-29727-NX66CDL	SW8270D	1-Methylnaphthalene	250000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	2-Methylnaphthalene	430000			DNR	<del></del>
Q4-15.0	08-29727-NX66CDL	SW8270D	Acenaphthene	300000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Acenaphthylene	36000		1 0	DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Anthracene	77000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Benzo(a)anthracene	110000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Benzo(a)pyrene	81000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Benzo(b)fluoranthene	73000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Benzo(g,h,i)perylene	36000		T U	DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Benzo(k)fluoranthene	51000		<del>  </del>	DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Chrysene	72000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Dibenz(a,h)anthracene	36000		U	DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Dibenzofuran	130000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Fluoranthene	320000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Fluorene	180000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Indeno(1,2,3-cd)pyrene	36000		<del>  U</del>	DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Phenanthrene	650000			DNR	11
Q4-15.0	08-29727-NX66CDL	SW8270D	Pyrene	300000		<del></del>	DNR	11
Q5-18.0	08-29733-NX66I	SW8270D	Phenanthrene		ug/kg	E	DNR	20
Q5-18.0	08-29733-NX66IDL	SW8270D	1-Methylnaphthalene		ug/kg		DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	2-Methylnaphthalene		ug/kg		DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Acenaphthene	2600		1	DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Acenaphthylene		ug/kg	U	DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Anthracene		ug/kg		DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Benzo(a)anthracene		ug/kg		DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Benzo(a)pyrene		ug/kg		DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Benzo(b)fluoranthene		ug/kg	1	DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Benzo(g,h,i)perylene		ug/kg		DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Benzo(k)fluoranthene		ug/kg		DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Chrysene		ug/kg		DNR	11

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	RC
Q5-18.0	08-29733-NX66IDL	SW8270D	Dibenz(a,h)anthracene		ug/kg	U	DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Dibenzofuran		ug/kg		DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Fluoranthene		ug/kg	ļ	DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Fluorene		ug/kg		DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg	U	DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Naphthalene		ug/kg		DNR	11
Q5-18.0	08-29733-NX66IDL	SW8270D	Pyrene		ug/kg	<u> </u>	DNR	11
Q6-4.0	08-29736-NX66K	SW8270D	Acenaphthene	770000		E	DNR	20
Q6-4.0	08-29736-NX66K	SW8270D	Anthracene	1600000		E	DNR	20
Q6-4.0	08-29736-NX66K	SW8270D	Benzo(a)anthracene	1600000		E	DNR	20
Q6-4.0	08-29736-NX66K	SW8270D	Benzo(a)pyrene	1800000		E	DNR	20
Q6-4.0	08-29736-NX66K	SW8270D	Benzo(b)fluoranthene	1400000		E	DNR	20
Q6-4.0	08-29736-NX66K	SW8270D	Benzo(g,h,i)perylene	820000		E	DNR	20
Q6-4.0	08-29736-NX66K	SW8270D	Benzo(k)fluoranthene	1600000		E	DNR	20
Q6-4.0	08-29736-NX66K	SW8270D	Chrysene	1900000		E	DNR	20
Q6-4.0	08-29736-NX66K	SW8270D	Fluoranthene	2400000		E	DNR	20
Q6-4.0	08-29736-NX66K	SW8270D	Phenanthrene	2800000		E	DNR	20
Q6-4.0	08-29736-NX66K	SW8270D	Pyrene	3400000		E	DNR	20
Q6-4.0	08-29736-NX66KDL	SW8270D	1-Methylnaphthalene		ug/kg	U	DNR	11
Q6-4.0	08-29736-NX66KDL	SW8270D	2-Methylnaphthalene		ug/kg		DNR	11
Q6-4.0	08-29736-NX66KDL	SW8270D	Acenaphthylene	92000		<u> </u>	DNR	11
Q6-4.0	08-29736-NX66KDL	SW8270D	Dibenz(a,h)anthracene	190000			DNR	11
Q6-4.0	08-29736-NX66KDL	SW8270D	Dibenzofuran	92000		U	DNR	11
Q6-4.0	08-29736-NX66KDL	SW8270D	Fluorene	170000			DNR	11
Q6-4.0	08-29736-NX66KDL	SW8270D	Indeno(1,2,3-cd)pyrene	840000			DNR	11
Q6-4.0	08-29736-NX66KDL	SW8270D	Naphthalene	92000		U	DNR	11
Q7-4.0	08-29740-NX66O	SW8270D	Benzo(a)anthracene	22000		Е	DNR	20
Q7-4.0	08-29740-NX66O	SW8270D	Benzo(a)pyrene	22000		E	DNR	20
Q7-4.0	08-29740-NX66O	SW8270D	Benzo(b)fluoranthene	21000		E	DNR	20
Q7-4.0	08-29740-NX66O	SW8270D	Benzo(k)fluoranthene	19000		E	DNR	20
Q7-4.0	08-29740-NX66O	SW8270D	Chrysene	23000		E	DNR	20
Q7-4.0	08-29740-NX66O	SW8270D	Fluoranthene	28000		E	DNR	20
Q7-4.0	08-29740-NX66O	SW8270D	Phenanthrene	29000		E	DNR	20
Q7-4.0	08-29740-NX66O	SW8270D	Pyrene	37000		Е	DNR	20
Q7-4.0	08-29740-NX66ODL	SW8270D	1-Methylnaphthalene		ug/kg	<u> </u>	DNR	11
Q7-4.0	08-29740-NX66ODL	SW8270D	2-Methylnaphthalene		ug/kg		DNR	11
Q7-4.0	08-29740-NX66ODL	SW8270D	Acenaphthene		ug/kg		DNR	11
Q7-4.0	08-29740-NX66ODL	SW8270D	Acenaphthylene		ug/kg	U	DNR	11
Q7-4.0	08-29740-NX66ODL	SW8270D	Anthracene		ug/kg		DNR	11
Q7-4.0	08-29740-NX66ODL	SW8270D	Benzo(g,h,i)perylene	17000			DNR	11
Q7-4.0	08-29740-NX66ODL	SW8270D	Dibenz(a,h)anthracene		ug/kg		DNR	11
Q7-4.0	08-29740-NX66ODL	SW8270D	Dibenzofuran		ug/kg		DNR	11
Q7-4.0	08-29740-NX66ODL	SW8270D	Fluorene		ug/kg		DNR	11
Q7-4.0	08-29740-NX66ODL	SW8270D	Indeno(1,2,3-cd)pyrene	13000			DNR	11
Q7-4.0	08-29740-NX66ODL	SW8270D	Naphthalene		ug/kg		DNR	11
Q7-5.5	08-29741-NX66P	SW8270D	Naphthalene	290000		E	DNR	20
Q7-5.5	08-29741-NX66P	SW8270D	Phenanthrene	240000		E	DNR	20
Q7-5.5	08-29741-NX66P	SW8270D	Pyrene	190000		<u>E</u>	DNR	20
Q7-5.5	08-29741-NX66PDL	SW8270D	1-Methylnaphthalene	59000			DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	2-Methylnaphthalene	84000			DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Acenaphthene	65000		ļ	DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Acenaphthylene	11000	ug/kg	U	DNR	11

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	RC
Q7-5.5	08-29741-NX66PDL	SW8270D	Anthracene	61000	ug/kg		DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Benzo(a)anthracene	86000	ug/kg		DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Benzo(a)pyrene	110000			DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Benzo(b)fluoranthene	57000	ug/kg		DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Benzo(g,h,i)perylene	72000	ug/kg		DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Benzo(k)fluoranthene	40000	ug/kg		DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Chrysene	110000			DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Dibenz(a,h)anthracene		ug/kg		DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Dibenzofuran	11000	ug/kg	U	DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Fluoranthene	140000	ug/kg		DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Fluorene		ug/kg		DNR	11
Q7-5.5	08-29741-NX66PDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg		DNR	11
Q9-28.0	08-29772-NX71F	SW8270D	Naphthalene	5200	ug/kg	E	DNR	20
Q9-28.0	08-29772-NX71FDL	SW8270D	1-Methylnaphthalene	520	ug/kg		DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	2-Methylnaphthalene		ug/kg		DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Acenaphthene	180	ug/kg		DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Acenaphthylene	180	ug/kg	Ū	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Anthracene		ug/kg	U	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Benzo(a)anthracene		ug/kg	U	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Benzo(a)pyrene		ug/kg	U	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Benzo(b)fluoranthene		ug/kg	U	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Benzo(g,h,i)perylene		ug/kg	TU	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Benzo(k)fluoranthene		ug/kg	U	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Chrysene		ug/kg	U	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Dibenz(a,h)anthracene		ug/kg	1 0	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Dibenzofuran		ug/kg	Ū	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Fluoranthene		ug/kg	U	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Fluorene		ug/kg	U	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg	U	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Phenanthrene		ug/kg	U	DNR	11
Q9-28.0	08-29772-NX71FDL	SW8270D	Pyrene		ug/kg	U	DNR	11
Q10-5.0	08-29773-NX71G	SW8270D	Anthracene		ug/kg	E	DNR	20
Q10-5.0	08-29773-NX71G	SW8270D	Benzo(a)anthracene	16000	ug/kg	ES	DNR	20
Q10-5.0	08-29773-NX71G	SW8270D	Benzo(a)pyrene	16000	ug/kg	ES	DNR	20
Q10-5.0	08-29773-NX71G	SW8270D	Benzo(b)fluoranthene	16000	ug/kg	ES	DNR	20
Q10-5.0	08-29773-NX71G	SW8270D	Benzo(k)fluoranthene	34000		ES	DNR	20
Q10-5.0	08-29773-NX71G	SW8270D	Chrysene	15000		ES	DNR	20
Q10-5.0	08-29773-NX71G	SW8270D	Fluoranthene	17000		ES	DNR	20
Q10-5.0	08-29773-NX71G	SW8270D	Indeno(1,2,3-cd)pyrene		ug/kg	E	DNR	20
Q10-5.0	08-29773-NX71G	SW8270D	Phenanthrene	14000		Е	DNR	20
Q10-5.0	08-29773-NX71G	SW8270D	Pyrene	19000	ug/kg	ES	DNR	20
Q10-5.0	08-29773-NX71GDL	SW8270D	1-Methylnaphthalene	2000	ug/kg		DNR	11
Q10-5.0	08-29773-NX71GDL	SW8270D	2-Methylnaphthalene		ug/kg		DNR	11
Q10-5.0	08-29773-NX71GDL	SW8270D	Acenaphthene		ug/kg		DNR	11
Q10-5.0	08-29773-NX71GDL	SW8270D	Acenaphthylene		ug/kg	U	DNR	11
Q10-5.0	08-29773-NX71GDL	SW8270D	Benzo(g,h,i)perylene	12000		T	DNR	11
Q10-5.0	08-29773-NX71GDL	SW8270D	Dibenz(a,h)anthracene		ug/kg	1	DNR	11
Q10-5.0	08-29773-NX71GDL	SW8270D	Dibenzofuran		ug/kg		DNR	11
Q10-5.0	08-29773-NX71GDL	SW8270D	Fluorene		ug/kg	1	DNR	11
Q10-5.0	08-29773-NX71GDL	SW8270D	Naphthalene		ug/kg		DNR	11
Q14-6.5	08-29837-NX79D	SW8270D	Phenanthrene	63000		Ē	DNR	20
Q14-6.5	08-29837-NX79D	SW8270D	Pyrene	62000		E	DNR	20

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	RC
Q14-6.5	08-29837-NX79DDL	SW8270D	1-Methylnaphthalene		ug/kg		DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	2-Methylnaphthalene		ug/kg		DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Acenaphthene		ug/kg		DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Acenaphthylene		ug/kg	U	DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Anthracene		ug/kg		DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Benzo(a)anthracene	25000			DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Benzo(a)pyrene	31000			DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Benzo(b)fluoranthene		ug/kg		DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Benzo(g,h,i)perylene	20000			DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Benzo(k)fluoranthene		ug/kg		DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Chrysene	33000			DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Dibenz(a,h)anthracene		ug/kg		DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Dibenzofuran		ug/kg	U	DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Fluoranthene	44000			DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Fluorene		ug/kg		DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Indeno(1,2,3-cd)pyrene	12000	ug/kg		DNR	11
Q14-6.5	08-29837-NX79DDL	SW8270D	Naphthalene	34000			DNR	11
Q15-4.0	08-29840-NX79G	SW8270D	Phenanthrene	60000	ug/kg	E	DNR	20
Q15-4.0	08-29840-NX79G	SW8270D	Pyrene	76000		E	DNR	20
Q15-4.0	08-29840-NX79GDL	SW8270D	1-Methylnaphthalene	7600	ug/kg		DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	2-Methylnaphthalene		ug/kg		DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Acenaphthene	8900	ug/kg		DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Acenaphthylene	2400	ug/kg	U	DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Anthracene	13000			DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Benzo(a)anthracene	34000			DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Benzo(a)pyrene	44000	ug/kg		DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Benzo(b)fluoranthene	27000	ug/kg		DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Benzo(g,h,i)perylene	28000			DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Benzo(k)fluoranthene	26000			DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Chrysene	50000			DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Dibenz(a,h)anthracene		ug/kg		DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Dibenzofuran		ug/kg	U	DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Fluoranthene	54000			DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Fluorene		ug/kg		DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Indeno(1,2,3-cd)pyrene	18000	ug/kg		DNR	11
Q15-4.0	08-29840-NX79GDL	SW8270D	Naphthalene	14000			DNR	11
SD-3	08-29852 <b>-</b> NX79S	SW8021B	Benzene		ug/kg		J	9
SD-3	08-29852-NX79S	SW8270D	1-Methylnaphthalene	220000	ug/kg	E	DNR	20
SD-3	08-29852-NX79S	SW8270D	2-Methylnaphthalene	310000		E	DNR	20
SD-3	08-29852-NX79S	SW8270D	Acenaphthene	190000		E	DNR	20
SD-3	08-29852-NX79S	SW8270D	Dibenzofuran	76000			J	9
SD-3	08-29852-NX79S	SW8270D	Fluoranthene	200000		E	DNR	20
SD-3	08-29852-NX79S	SW8270D	Fluorene	140000		E	DNR	20
SD-3	08-29852-NX79S	SW8270D	Naphthalene	480000		E	DNR	20
SD-3	08-29852-NX79S	SW8270D	Phenanthrene	360000		E	DNR	20
SD-3	08-29852-NX79S	SW8270D	Pyrene	170000		E	DNR	20
SD-3	08-29852-NX79SDL	SW8270D	Acenaphthylene	12000		U	DNR	11
SD-3	08-29852-NX79SDL	SW8270D	Anthracene	70000	ug/kg		DNR	11
SD-3	08-29852-NX79SDL	SW8270D	Benzo(a)anthracene	68000			DNR	11
SD-3	08-29852-NX79SDL	SW8270D	Benzo(a)pyrene	48000			DNR	11
SD-3	08-29852-NX79SDL	SW8270D	Benzo(b)fluoranthene	33000			DNR	11
SD-3	08-29852-NX79SDL	SW8270D	Benzo(g,h,i)perylene	16000	ug/kg		DNR	11

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Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	RC
SD-3	08-29852-NX79SDL	SW8270D	Benzo(k)fluoranthene	50000	ug/kg		DNR	11
SD-3	08-29852-NX79SDL	SW8270D	Chrysene	49000	ug/kg		DNR	11
SD-3	08-29852-NX79SDL	SW8270D	Dibenz(a,h)anthracene	12000	ug/kg	U	DNR	11
SD-3	08-29852-NX79SDL	SW8270D	Dibenzofuran	72000	ug/kg		DNR	11
SD-3	08-29852-NX79SDL	SW8270D	Indeno(1,2,3-cd)pyrene	15000	ug/kg		DNR	11
Q4-W	08-29958-NY02K	SW6010B	Arsenic	0.05	mg/l	U	DNR	11
Q4-W	08-29958-NY02K	SW6010B	Lead		? mg/l	U	DNR	11
Q4-W	08-29958-NY02K	SW8270D	1-Methylnaphthalene	640	ug/L	ES	DNR	10
Q4-W	08-29958-NY02K	SW8270D	2-Methylnaphthalene		ug/L	ES	DNR	10
Q4-W	08-29958-NY02K	SW8270D	Acenaphthene	200	ug/L	E	DNR	10
Q4-W	08-29958-NY02K	SW8270D	Acenaphthylene		ug/L		DNR	10
Q4-W	08-29958-NY02K	SW8270D	Anthracene	45	ug/L		DNR	10
Q4-W	08-29958-NY02K	SW8270D	Benzo(a)anthracene		ug/L		DNR	10
Q4-W	08-29958-NY02K	SW8270D	Benzo(a)pyrene	21	ug/L		DNR	10
Q4-W	08-29958-NY02K	SW8270D	Benzo(b)fluoranthene	20	ug/L		DNR	10
Q4-W	08-29958-NY02K	SW8270D	Benzo(g,h,i)perylene	7	ug/L		DNR	10
Q4-W	08-29958-NY02K	SW8270D	Benzo(k)fluoranthene	14	ug/L		DNR	10
Q4-W	08-29958-NY02K	SW8270D	Chrysene		ug/L		DNR	10
Q4-W	08-29958-NY02K	SW8270D	Dibenz(a,h)anthracene		ug/L		DNR	10
Q4-W	08-29958-NY02K	SW8270D	Dibenzofuran	110	ug/L	E	DNR	10
Q4-W	08-29958-NY02K	SW8270D	Fluoranthene		ug/L	E	DNR	10
Q4-W	08-29958-NY02K	SW8270D	Fluorene		ug/L	E	DNR	10
Q4-W	08-29958-NY02K	SW8270D	Indeno(1,2,3-cd)pyrene	7	ug/L		DNR	10
Q4-W	08-29958-NY02K	SW8270D	Naphthalene	87000		ES	DNR	10
Q4-W	08-29958-NY02K	SW8270D	Phenanthrene	220	ug/L	E	DNR	10
Q4-W	08-29958-NY02K	SW8270D	Pyrene	79	ug/L		DNR	10
Q4-W	08-29958-NY02KDL	NWTPH-Dx	Diesel Range Hydrocarbons		mg/L		DNR	11
Q4-W	08-29958-NY02KDL	NWTPH-Dx	Motor Oil	2.5	mg/L	U	DNR	11
Q4-W	08-29958-NY02KRE	SW8270D	1-Methylnaphthalene	710	ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	2-Methylnaphthalene	1100	ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Acenaphthene	620	ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Acenaphthylene	20	ug/L	U	UJ	1
Q4-W	08-29958-NY02KRE	SW8270D	Anthracene		ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Benzo(a)anthracene		ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Benzo(a)pyrene	65	ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Benzo(b)fluoranthene	60	ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Benzo(g,h,i)perylene	28	ug/L		Ĵ	1
Q4-W	08-29958-NY02KRE	SW8270D	Benzo(k)fluoranthene	36	ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Chrysene	60	ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Dibenz(a,h)anthracene	20	ug/L	U	UJ	1
Q4-W	08-29958-NY02KRE	SW8270D	Dibenzofuran	260	ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Fluoranthene		ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Fluorene	330	ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Indeno(1,2,3-cd)pyrene		ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Naphthalene	8300	ug/L	Е	DNR	20
Q4-W	08-29958-NY02KRE	SW8270D	Phenanthrene	840	ug/L		J	1
Q4-W	08-29958-NY02KRE	SW8270D	Pyrene	270	ug/L		J	1
Q4-W	08-29958-NY02KREDL	SW8270D	1-Methylnaphthalene	770	ug/L		DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	2-Methylnaphthalene	1200	ug/L		DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Acenaphthene	690	ug/L		DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Acenaphthylene	200	ug/L	U	DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Anthracene	200	ug/L	U	DNR	11

Γ	T	1	Quendan Terminai - Bor	1	T	Lab	T	DV Qual
Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	RC
Q4-W	08-29958-NY02KREDL	SW8270D	Benzo(a)anthracene		ug/L	U	DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Benzo(a)pyrene		ug/L	U	DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Benzo(b)fluoranthene		ug/L	U	DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Benzo(g,h,i)perylene		ug/L	U	DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Benzo(k)fluoranthene		ug/L	U	DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Chrysene	200	ug/L	TU	DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Dibenz(a,h)anthracene	200	ug/L	U	DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Dibenzofuran		ug/L		DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Fluoranthene		ug/L		DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Fluorene		ug/L		DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/L	U	DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Naphthalene		ug/L		J	1
Q4-W	08-29958-NY02KREDL	SW8270D	Phenanthrene		ug/L		DNR	11
Q4-W	08-29958-NY02KREDL	SW8270D	Pyrene		ug/L		DNR	11
Q9-W	08-29959-NY02L	SW8270D	1-Methylnaphthalene		ug/L	E	DNR	20
Q9-W	08-29959-NY02L	SW8270D	2-Methylnaphthalene	12000		E	DNR	20
Q9-W	08-29959-NY02L	SW8270D	Acenaphthene	6700		E	DNR	20
Q9-W	08-29959-NY02L	SW8270D	Acenaphthylene		ug/L		J	10
Q9-W	08-29959-NY02L	SW8270D	Anthracene		ug/L		J	10
Q9-W	08-29959-NY02L	SW8270D	Benzo(a)anthracene	3100	ua/L	<del>                                     </del>	J	10
Q9-W	08-29959-NY02L	SW8270D	Benzo(a)pyrene	2000	ug/L		J	10
Q9-W	08-29959-NY02L	SW8270D	Benzo(b)fluoranthene	1900	ug/L		J	10
Q9-W	08-29959-NY02L	SW8270D	Benzo(g,h,i)perylene		ug/L		J	10
Q9-W	08-29959-NY02L	SW8270D	Benzo(k)fluoranthene	1400			J	10
Q9-W	08-29959-NY02L	SW8270D	Chrysene	2200			J	10
Q9-W	08-29959-NY02L	SW8270D	Dibenz(a,h)anthracene		ug/L		J	10
Q9-W	08-29959-NY02L	SW8270D	Dibenzofuran	3800			J	10
Q9-W	08-29959-NY02L	SW8270D	Fluoranthene	7500		E	DNR	20
Q9-W	08-29959-NY02L	SW8270D	Fluorene	11000		E	DNR	20
Q9-W	08-29959-NY02L	SW8270D	Indeno(1,2,3-cd)pyrene		ug/L	<del>                                     </del>	J	10
Q9-W	08-29959-NY02L	SW8270D	Naphthalene	120000		ES	DNR	20
Q9-W	08-29959-NY02L	SW8270D	Phenanthrene	14000		E	DNR	20
Q9-W	08-29959-NY02L	SW8270D	Pyrene	6200		E	DNR	20
Q9-W	08-29959-NY02LDL	SW8270D	1-Methylnaphthalene	10000			J	10
Q9-W	08-29959-NY02LDL	SW8270D	2-Methylnaphthalene	17000			J	10
Q9-W	08-29959-NY02LDL	SW8270D	Acenaphthene	10000			J	10
Q9-W	08-29959-NY02LDL	SW8270D	Acenaphthylene	1100			DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Anthracene	2800			DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Benzo(a)anthracene	3500			DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Benzo(a)pyrene	2300			DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Benzo(b)fluoranthene	1800			DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Benzo(g,h,i)perylene	1000			DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Benzo(k)fluoranthene	1500		1	DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Chrysene	2400			DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Dibenz(a,h)anthracene		ug/L	U	DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Dibenzofuran	5000			DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Fluoranthene	10000			J	10
Q9-W	08-29959-NY02LDL	SW8270D	Fluorene	7300			J	10
Q9-W	08-29959-NY02LDL	SW8270D	Indeno(1,2,3-cd)pyrene		ug/L		DNR	11
Q9-W	08-29959-NY02LDL	SW8270D	Naphthalene	65000		Е	DNR	20
Q9-W	08-29959-NY02LDL	SW8270D	Phenanthrene	23000			J	10
Q9-W	08-29959-NY02LDL	SW8270D	Pyrene	11000	ug/L		J	10

Sample ID		<u> </u>					Lab		DV Qual
Q9-W   Q8-29959-NYQ2LDL2   SW8270D   Asnaphthylene   19000lugit   DNR   11	Sample ID		Method	Analyte			Flag	DV Qual	RC
C9-W   08-2955-NY02LDL2   SW8270D   Acenaphthene   10000 log.t.   DNR   11	Q9-W	08-29959-NY02LDL2	SW8270D	1-Methylnaphthalene				DNR	11
G9-W   08-29959-NY02LDL2   SW8270D   Abenaphtylene   3000 lugit   U   DNR   11				2-Methylnaphthalene				DNR	
G9-W   G9-2965-NYOZLD.2   SW8270D   Benzo(g), primare   3000 lug/L   DNR   11				Acenaphthene					
Q9-W   Q8-29959-NYQ2,DL2   SW8270D   Benzo(a)anthracene   3400 light		08-29959-NY02LDL2					U		
G9-W   G8-29958-NY02LDL2   SW8270D   Benzo(a)pyrene   3000 lug/L   U DNR   11		08-29959-NY02LDL2		Anthracene			U	DNR	
Q3-W   Q8-29959-NYQ2LDL2   SW8270D   Benzo(gh.li)perylene   3000 lgg/L   U DNR   11		08-29959-NY02LDL2		Benzo(a)anthracene					
Q3-W   Q8-29959-NYQ2_DL2   SW8270D   Benzo(gh_liperylene   3000 lug/L   U   DNR   11	Q9-W	08-29959-NY02LDL2	SW8270D	Benzo(a)pyrene	3000	ug/L	U	DNR	
C9-W         08-29958-NYOZLDL2         SW8270D         Berazok/jituoranthene         3000 lggl.         U         DNR         11           C9-W         08-29959-NYOZLDL2         SW8270D         Dibenz(a,i)janthracene         3000 lggl.         U         DNR         11           C9-W         08-29959-NYOZLDL2         SW8270D         Dibenz(a,i)janthracene         3000 lggl.         U         DNR         11           C9-W         08-29959-NYOZLDL2         SW8270D         Fluoranthene         10000 lggl.         DNR         11           C9-W         08-29959-NYOZLDL2         SW8270D         Fluoranthene         3000 lggl.         U         DNR         11           C9-W         08-29959-NYOZLDL2         SW8270D         Naphthelane         45000 lggl.         J         10           C9-W         08-29959-NYOZLDL2         SW8270D         Naphthelane         45000 lggl.         DNR         11           C9-W         08-29959-NYOZLDL2         SW8270D         Pyrene         10000 lggl.         DNR         11           C12-W         08-29959-NYOZLDL2         SW8270D         Pyrene         10000 lggl.         DNR         11           C12-W         08-29959-NYOZLD         SW8270D         Pyrene         10000 lggl.				Benzo(b)fluoranthene			U		
Q9-W	Q9-W	08-29959-NY02LDL2	SW8270D	Benzo(g,h,i)perylene			U	DNR	
Q9-W		08-29959-NY02LDL2	SW8270D	Benzo(k)fluoranthene			U	DNR	
Q9-W     08-29959-NY02LDL2   SW8270D   Dibenzofuran     5100 log/L   DNR   11	Q9-W	08-29959-NY02LDL2	SW8270D				U	DNR	
Q9-W         08-29959-NYO2LDL2         SW8270D         Fluoranthene         10000 lggl.         DNR         11           Q9-W         08-29959-NYO2LDL2         SW8270D         Indence(1,2,3-cd)pyrene         6000 lugl.         U         DNR         11           Q9-W         08-29959-NYO2LDL2         SW8270D         Naphthalene         45000 lugl.         U         DNR         11           Q9-W         08-29959-NYO2LDL2         SW8270D         Naphthalene         45000 lugl.         DNR         11           Q9-W         08-29959-NYO2LDL2         SW8270D         Pyrene         10000 lugl.         DNR         11           Q12-W         08-29959-NYO2LDL2         SW8270D         Pyrene         10000 lugl.         DNR         11           Q12-W         08-29959-NYO2LD         SW8270D         Pyrene         10000 lugl.         DNR         11           Q12-W         08-29960-NYO2M         SW8010B         Arsenic         0.05 lmgl         U         DNR         11           Q12-W         08-29960-NYO2M         SW8270D         Acenaphthylene         3.2 lugl.         DNR         10           Q12-W         08-29960-NYO2M         SW8270D         Acenaphthylene         1 lugl.         U         DNR         1	Q9-W	08-29959-NY02LDL2	SW8270D	Dibenz(a,h)anthracene			U	DNR	11
Q9-W         08-29959-NY02LDL2         SW8270D         Fluorene         6000 lugl.         DNR         11           Q9-W         08-29959-NY02LDL2         SW8270D         Indeno(1,2,3-cd)pyrene         3000 lugl.         U         DNR         11           Q9-W         08-29959-NY02LDL2         SW8270D         Naphthalene         45000 lugl.         DNR         11           Q3-W         08-29959-NY02LDL2         SW8270D         Pyrene         10000 lugl.         DNR         11           Q3-W         08-29950-NY02LDL2         SW8270D         Pyrene         10000 lugl.         DNR         11           Q3-W         08-29950-NY02M         SW8010B         Arsenic         0.05 lmgl.         U         DNR         11           Q12-W         08-29960-NY02M         SW8010B         Lead         0.02 lngl.         U         DNR         11           Q12-W         08-29960-NY02M         SW8270D         1-Methylnaphthalene         2.1 lugl.         DNR         10           Q12-W         08-29960-NY02M         SW8270D         2-Methylnaphthalene         4.2 lugl.         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Acenaphthylene         1 lugl.         U         DNR         10<	Q9-W	08-29959-NY02LDL2	SW8270D	Dibenzofuran	5100	ug/L		DNR	11
G9-W   08-29959-NY02LDL2   SW8270D   Indeno(1,2,3-cd)pyrene   3000 lug/L   J   J   10	Q9-W	08-29959-NY02LDL2	SW8270D	Fluoranthene				DNR	11
Q9-W         08-29959-NYO2LDL2         SW8270D         Naphthalene         45000 lug/L         J         10           Q9-W         08-29959-NYO2LDL2         SW8270D         Phenanthrene         26000 lug/L         DNR         11           Q9-W         08-29950-NY02M         SW8270D         Pyrene         10000 lug/L         DNR         11           Q12-W         08-29960-NY02M         SW8270D         Nebroine         0.05 mg/l         U         DNR         11           Q12-W         08-29960-NY02M         SW8270D         1-Melthylnaphthalene         2.1 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         2-Methylnaphthalene         3.2 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Acenaphthylene         4.2 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Acenaphthylene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(a)anthracene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(b)Hiuoranthene         1 ug/L         U         DNR <td< td=""><td>Q9-W</td><td>08-29959-NY02LDL2</td><td>SW8270D</td><td>Fluorene</td><td>6000</td><td>ug/L</td><td></td><td>DNR</td><td>11</td></td<>	Q9-W	08-29959-NY02LDL2	SW8270D	Fluorene	6000	ug/L		DNR	11
Q9-W   Q8-29958-NYO2LDL2	Q9-W	08-29959-NY02LDL2	SW8270D	Indeno(1,2,3-cd)pyrene	3000	ug/L	U	DNR	11
Q9-W     Q8-29958-NY02LDL2   SW8270D   Pyrene   10000 lg/L   DNR   11   Q12-W   Q8-29960-NY02M   SW6010B   Arsenic   Q.02 mg/l   U DNR   11   Q12-W   Q8-29960-NY02M   SW6270D   1-Methylnaphthalene   Q.1 lg/L   DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   2-Methylnaphthalene   Q.1 lg/L   DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Acenaphthene   Q.1 lg/L   DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Acenaphthene   Q.1 lg/L   U DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Acenaphthene   Q.1 lg/L   U DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Acenaphthalene   Q1-lg/L   U DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Acenaphthalene   Q1-lg/L   U DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Benzo(a)pyrene   Q1-lg/L   U DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Benzo(a)pyrene   Q1-lg/L   U DNR   Q1-lg-L   Q8-29960-NY02M   SW8270D   Benzo(a)pyrene   Q1-lg/L   U DNR   Q1-lg-L   Q8-29960-NY02M   SW8270D   Benzo(b)lororanthene   Q1-lg/L   U DNR   Q1-lg-L   Q1-	Q9-W	08-29959-NY02LDL2	SW8270D	Naphthalene	45000	ug/L		J	10
Q9-W     Q8-29958-NY02LDL2   SW8270D   Pyrene   10000 lg/L   DNR   11   Q12-W   Q8-29960-NY02M   SW6010B   Arsenic   Q.02 mg/l   U DNR   11   Q12-W   Q8-29960-NY02M   SW6270D   1-Methylnaphthalene   Q.1 lg/L   DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   2-Methylnaphthalene   Q.1 lg/L   DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Acenaphthene   Q.1 lg/L   DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Acenaphthene   Q.1 lg/L   U DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Acenaphthene   Q.1 lg/L   U DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Acenaphthalene   Q1-lg/L   U DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Acenaphthalene   Q1-lg/L   U DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Benzo(a)pyrene   Q1-lg/L   U DNR   10   Q12-W   Q8-29960-NY02M   SW8270D   Benzo(a)pyrene   Q1-lg/L   U DNR   Q1-lg-L   Q8-29960-NY02M   SW8270D   Benzo(a)pyrene   Q1-lg/L   U DNR   Q1-lg-L   Q8-29960-NY02M   SW8270D   Benzo(b)lororanthene   Q1-lg/L   U DNR   Q1-lg-L   Q1-	Q9-W	08-29959-NY02LDL2	SW8270D	Phenanthrene	26000	ug/L		DNR	11
Q12-W   08-29960-NY02M   SW8010B   Arsenic   0.05 mg/l   U   DNR   11	Q9-W	08-29959-NY02LDL2	SW8270D	Pyrene				DNR	11
Q12-W   Q8-2996Q-NYQ2M   SW827DD   Lead   Q.02 mg/l   Q. DNR   11	Q12-W	08-29960-NY02M	SW6010B		0.05	mg/l	U	DNR	11
Q12-W   08-29960-NY02M   SW8270D   1-Methylnaphthalene   2.1 ug/L   DNR   10	Q12-W	08-29960-NY02M	SW6010B	Lead			U	DNR	11
Q12-W         08-29960-NY02M         SW8270D         2-Methylnaphthalene         3.2 lug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Acenaphthene         4.2 lug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Acenaphthylene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Anthracene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(a)anthracene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dienzo(s)fluoranthene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenzo(s)fl		08-29960-NY02M	SW8270D	1-Methylnaphthalene	2.1	ug/L		DNR	10
Q12-W         08-29960-NY02M         SW8270D         Acenaphthene         4.2 lg/fL         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Acenaphthylene         1 lg/fL         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Anthracene         1 lg/fL         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(a)anthracene         1 lug/fL         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(g)pyrene         1 lug/fL         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(g)hijperylene         1 lug/fL         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(g,hijperylene         1 lug/fL         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(g,hijperylene         1 lug/fL         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(g,hijperylene         1 lug/fL         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D		08-29960-NY02M	SW8270D					DNR	10
Q12-W         08-29960-NY02M         SW8270D         Acenaphthylene         1 lug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Anthracene         1 lug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(a)anthracene         1 lug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(a)pyrene         1 lug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(g)hijoeranthene         1 lug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(k)fluoranthene         1 lug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenz(k)fluoranthene         1 lug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenz(a,h)anthracene         1 lug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenzofuran         1 lug/L         U         DNR         10           Q12-W         08-29960-NY02M <td< td=""><td></td><td>08-29960-NY02M</td><td>SW8270D</td><td>Acenaphthene</td><td></td><td></td><td></td><td>DNR</td><td>10</td></td<>		08-29960-NY02M	SW8270D	Acenaphthene				DNR	10
Q12-W         08-29960-NY02M         SW8270D         Anthracene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(a)anthracene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(a)pyrene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenzofuran         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Horere		08-29960-NY02M	SW8270D	Acenaphthylene	1		U	DNR	10
Q12-W   Q8-29960-NY02M   SW8270D   Benzo(a)anthracene   1 ug/L   U DNR   10		08-29960-NY02M	SW8270D		1		Ū	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Benzo(a)pyrene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(g,hiperplene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(g,hiperplene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Chrysene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenz(a,h)anthracene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenzofuran         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.1 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Indeno(1,2,3-cd)pyrene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Naphthalene         17		08-29960-NY02M	SW8270D	Benzo(a)anthracene			U	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Benzo(g,h,i)perylene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Chrysene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenz(a,h)anthracene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenzofuran         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.1 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Plenarthrene         3.1 ug/L         U         DNR         10           Q12-W         08-29960-NY02MR         SW8270D         Pyrene         1 ug/L         DNR </td <td></td> <td>08-29960-NY02M</td> <td>SW8270D</td> <td></td> <td>1</td> <td>ug/L</td> <td>U</td> <td>DNR</td> <td>10</td>		08-29960-NY02M	SW8270D		1	ug/L	U	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Benzo(g,h,i)perylene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Chrysene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenz(a,h)anthracene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenzofuran         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.1 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Plenarthrene         3.1 ug/L         U         DNR         10           Q12-W         08-29960-NY02MR         SW8270D         Pyrene         1 ug/L         DNR </td <td>Q12-W</td> <td>08-29960-NY02M</td> <td>SW8270D</td> <td>Benzo(b)fluoranthene</td> <td>1</td> <td>ug/L</td> <td>U</td> <td>DNR</td> <td>10</td>	Q12-W	08-29960-NY02M	SW8270D	Benzo(b)fluoranthene	1	ug/L	U	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Chrysene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenz(a,h)anthracene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenzofuran         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.1 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluorene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Indeno(1,2,3-cd)pyrene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Naphthalene         17 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Pyrene         1 ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         Pyrene         1 ug/L         J         1		08-29960-NY02M	SW8270D	Benzo(g,h,i)perylene			Ū	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Chrysene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenz(a,h)anthracene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluorene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluorene         1.5 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Indeno(1,2,3-cd)pyrene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Naphthalene         17 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Phenanthrene         3.1 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Pyrene         1 ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthalene         1.7 ug/L         J         1,9     <	Q12-W	08-29960-NY02M	SW8270D	Benzo(k)fluoranthene	1		U	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Dibenz(a,h)anthracene         1 ug/L         U DNR         10           Q12-W         08-29960-NY02M         SW8270D         Dibenzofuran         1 ug/L         U DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.1 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluorene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluorene         1 ug/L         U DNR         10           Q12-W         08-29960-NY02M         SW8270D         Naphthalene         17 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Phenanthrene         3.1 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Pyrene         1 ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         1-Methylnaphthalene         1.3 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         2-Methylnaphthalene         1.7 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8		08-29960-NY02M	SW8270D	Chrysene	1	ug/L	U	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Dibenzofuran         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.1 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluorene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Indeno(1,2,3-cd)pyrene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Naphthalene         17 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Phenanthrene         3.1 ug/L         DNR         10           Q12-W         08-29960-NY02MR         SW8270D         Pyrene         1 ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         Pyrene         1.3 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         2-Methylnaphthalene         1.7 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1 ug/L         U         U         J         1,9         1 </td <td></td> <td>08-29960-NY02M</td> <td>SW8270D</td> <td></td> <td></td> <td></td> <td>U</td> <td>DNR</td> <td>10</td>		08-29960-NY02M	SW8270D				U	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Fluoranthene         1.1         ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Fluorene         1.5         ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Indeno(1,2,3-cd)pyrene         1         ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Naphthalene         17         ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Phenanthrene         3.1         ug/L         DNR         10           Q12-W         08-29960-NY02MR         SW8270D         Pyrene         1         ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         1-Methylnaphthalene         1.3         ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthene         3.7         ug/L         J         1,9           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D </td <td></td> <td>08-29960-NY02M</td> <td>SW8270D</td> <td>Dibenzofuran</td> <td></td> <td></td> <td>U</td> <td>DNR</td> <td>10</td>		08-29960-NY02M	SW8270D	Dibenzofuran			U	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Fluorene         1.5 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Indeno(1,2,3-cd)pyrene         1 ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Naphthalene         17 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Phenanthrene         3.1 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Pyrene         1 ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         1-Methylnaphthalene         1.3 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         2-Methylnaphthalene         1.7 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthene         3.7 ug/L         J         1,9           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)anthracene         1 ug/L         U         UJ         1		08-29960-NY02M	SW8270D	Fluoranthene				DNR	10
Q12-W         08-29960-NY02M         SW8270D         Indeno(1,2,3-cd)pyrene         1         ug/L         U         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Naphthalene         17         ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Phenanthrene         3.1         ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         Pyrene         1         ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         1-Methylnaphthalene         1.3         ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthene         3.7         ug/L         J         1,3           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1         ug/L         U         UJ         1,3           Q12-W         08-29960-NY02MRE         SW8270D         Anthracene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)anthracene         1         ug/L         U         UJ         1           Q12-W         0		08-29960-NY02M		Fluorene			1	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Naphthalene         17 ug/L         DNR         10           Q12-W         08-29960-NY02M         SW8270D         Phenanthrene         3.1 ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         Pyrene         1 ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         1-Methylnaphthalene         1.3 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         2-Methylnaphthalene         1.7 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthene         3.7 ug/L         J         1,9           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Anthracene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)pyrene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         UJ         1		08-29960-NY02M	SW8270D	Indeno(1,2,3-cd)pyrene			U	DNR	10
Q12-W         08-29960-NY02M         SW8270D         Phenanthrene         3.1         ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         Pyrene         1         ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         1-Methylnaphthalene         1.3         ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthene         3.7         ug/L         J         1,9           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Anthracene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)anthracene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)pyrene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1         ug/L         U         UJ         1           Q12-	Q12-W	08-29960-NY02M	SW8270D	Naphthalene	17	ug/L		DNR	10
Q12-W         08-29960-NY02M         SW8270D         Pyrene         1 ug/L         DNR         10           Q12-W         08-29960-NY02MRE         SW8270D         1-Methylnaphthalene         1.3 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         2-Methylnaphthalene         1.7 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthene         3.7 ug/L         U         UJ         1,9           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)anthracene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)pyrene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 u		08-29960-NY02M	SW8270D	Phenanthrene				DNR	10
Q12-W         08-29960-NY02MRE         SW8270D         1-Methylnaphthalene         1.3         ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         2-Methylnaphthalene         1.7         ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Anthracene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)anthracene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)pyrene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(g,h,i)perylene         1         ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1         ug/L         U         U		08-29960-NY02M	SW8270D	Pyrene				DNR	10
Q12-W         08-29960-NY02MRE         SW8270D         2-Methylnaphthalene         1.7 ug/L         J         1           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthene         3.7 ug/L         J         1,9           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)anthracene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)pyrene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(g,h,i)perylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Chrysene         1 ug/L         U         UJ         1		08-29960-NY02MRE	SW8270D	1-Methylnaphthalene				J	1
Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthene         3.7 ug/L         J         1,9           Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Anthracene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)pyrene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(g,h,i)perylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Chrysene         1 ug/L         U         UJ         1		08-29960-NY02MRE	SW8270D	2-Methylnaphthalene			1	J	1
Q12-W         08-29960-NY02MRE         SW8270D         Acenaphthylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Anthracene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)anthracene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)pyrene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(g,h,i)perylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Chrysene         1 ug/L         U         UJ         1		1							1,9
Q12-W         08-29960-NY02MRE         SW8270D         Anthracene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)anthracene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)pyrene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(g,h,i)perylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Chrysene         1 ug/L         U         UJ         1				<u> </u>			U		1
Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)anthracene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)pyrene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(g,h,i)perylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Chrysene         1 ug/L         U         UJ         1				· <del>  </del>			U		1
Q12-W         08-29960-NY02MRE         SW8270D         Benzo(a)pyrene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(g,h,i)perylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Chrysene         1 ug/L         U         UJ         1		<del></del>					U		1
Q12-W         08-29960-NY02MRE         SW8270D         Benzo(b)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(g,h,i)perylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Chrysene         1 ug/L         U         UJ         1				<del></del>			U	UJ	1
Q12-W         08-29960-NY02MRE         SW8270D         Benzo(g,h,i)perylene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Chrysene         1 ug/L         U         UJ         1							U		1
Q12-W         08-29960-NY02MRE         SW8270D         Benzo(k)fluoranthene         1 ug/L         U         UJ         1           Q12-W         08-29960-NY02MRE         SW8270D         Chrysene         1 ug/L         U         UJ         1		<u> </u>					U		1
Q12-W 08-29960-NY02MRE SW8270D Chrysene 1 ug/L U UJ 1				· <del></del>			U		1
				1			U		1
₩12-99  00-2000-1010  0990/C  0990/C   Dinch2(α₁)βαιπιασσία   Ilug/C   O   OJ   I	Q12-W	08-29960-NY02MRE	SW8270D	Dibenz(a,h)anthracene		ug/L	U	UJ	1

Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Lab Flag	DV Qual	DV Qual RC
Q12-W	08-29960-NY02MRE	SW8270D	Dibenzofuran	1	ug/L	U	UJ	1
Q12-W	08-29960-NY02MRE	SW8270D	Fluoranthene	1	ug/L	Ū	UJ	1
Q12-W	08-29960-NY02MRE	SW8270D	Fluorene	1.1	ug/L		J	1
Q12-W	08-29960-NY02MRE	SW8270D	Indeno(1,2,3-cd)pyrene	1	ug/L	U	UJ	1
Q12-W	08-29960-NY02MRE	SW8270D	Naphthalene	7.8	ug/L		J	1,9
Q12-W	08-29960-NY02MRE	SW8270D	Phenanthrene	1.5	ug/L		J	1
Q12-W	08-29960-NY02MRE	SW8270D	Pyrene	1	ug/L	U	UJ	1,9
Q14-W	08-29961-NY02N	SW6010B	Arsenic	0.05	mg/l	U	DNR	11
Q14-W	08-29961-NY02N	SW6010B	Lead	0.02	mg/l	T	DNR	11
Q14-W	08-29961-NY02N	SW8270D	1-Methylnaphthalene		ug/L		DNR	10
Q14-W	08-29961-NY02N	SW8270D	2-Methylnaphthalene	1.7	ug/L		DNR	10
Q14-W	08-29961-NY02N	SW8270D	Acenaphthene	1	ug/L	U	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Acenaphthylene	1	ug/L	U	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Anthracene	1	ug/L	U	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Benzo(a)anthracene	1	ug/L	U	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Benzo(a)pyrene		ug/L	U	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Benzo(b)fluoranthene		ug/L	U	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Benzo(g,h,i)perylene	1	ug/L	Ū	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Benzo(k)fluoranthene	1	ug/L	U	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Chrysene	1	ug/L	Ū	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Dibenz(a,h)anthracene	1	ug/L	U	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Dibenzofuran	1	ug/L	Ū	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Fluoranthene	1	ug/L	U	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Fluorene	1	ug/L	Ū	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Indeno(1,2,3-cd)pyrene	1	ug/L	U	DNR	10
Q14-W	08-29961-NY02N	SW8270D	Naphthalene		ug/L		DNR	10
Q14-W	08-29961-NY02N	SW8270D	Phenanthrene	2.1	ug/L		DNR	10
Q14-W	08-29961-NY02N	SW8270D	Pyrene		ug/L	U	DNR	10
Q14-W	08-29961-NY02NRE	SW8270D	1-Methylnaphthalene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	2-Methylnaphthalene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Acenaphthene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Acenaphthylene		ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Anthracene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Benzo(a)anthracene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Benzo(a)pyrene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Benzo(b)fluoranthene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Benzo(g,h,i)perylene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Benzo(k)fluoranthene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Chrysene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Dibenz(a,h)anthracene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Dibenzofuran	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Fluoranthene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Fluorene	1	ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Indeno(1,2,3-cd)pyrene		ug/L	U	UJ	1
Q14-W	08-29961-NY02NRE	SW8270D	Naphthalene		ug/L		J	1
Q14-W	08-29961-NY02NRE	SW8270D	Phenanthrene		ug/L		J	1
Q14-W	08-29961-NY02NRE	SW8270D	Pyrene		ug/L		J	1
Q17-W	08-29962-NY02O	NWTPH-Dx	Diesel Range Hydrocarbons		mg/L	U	UJ	13
Q17-W	08-29962-NY02O	NWTPH-Dx	Motor Oil	0.5	mg/L	U	UJ	13
Q17-W	08-29962-NY02O	SW6010B	Arsenic	0.05		U	DNR	11
Q17-W	08-29962-NY02O	SW6010B	Lead	0.02		U	DNR	11
Q17-W	08-29962-NY02O	SW8270D	1-Methylnaphthalene		ug/L	U	DNR	10

						Lab		DV Qual
Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	RC
Q17-W	08-29962-NY02O	SW8270D	2-Methylnaphthalene	11	ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Acenaphthene		ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Acenaphthylene	1	ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Anthracene	1	ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Benzo(a)anthracene		ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Benzo(a)pyrene		ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Benzo(b)fluoranthene	1	ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Benzo(g,h,i)perylene	1	ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Benzo(k)fluoranthene	11	ug/L	U	DNR	10
Q17-W_	08-29962-NY02O	SW8270D	Chrysene	1 1	ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Dibenz(a,h)anthracene		ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Dibenzofuran		ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Fluoranthene		ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Fluorene		ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Indeno(1,2,3-cd)pyrene		ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Naphthalene		ug/L		DNR	10
Q17-W	08-29962-NY02O	SW8270D	Phenanthrene		ug/L	U	DNR	10
Q17-W	08-29962-NY02O	SW8270D	Pyrene		ug/L	U	DNR	10
Q17-W	08-29962-NY02ORE	NWTPH-Dx	Diesel Range Hydrocarbons		mg/L	U	DNR	11
Q17-W	08-29962-NY02ORE	NWTPH-Dx	Motor Oil	0.5	mg/L	U	DNR	11
Q17-W_	08-29962-NY02ORE	SW8270D	1-Methylnaphthalene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	2-Methylnaphthalene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Acenaphthene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Acenaphthylene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Anthracene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Benzo(a)anthracene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Benzo(a)pyrene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Benzo(b)fluoranthene		ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Benzo(g,h,i)perylene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Benzo(k)fluoranthene		ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Chrysene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Dibenz(a,h)anthracene		ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Dibenzofuran	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Fluoranthene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Fluorene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Indeno(1,2,3-cd)pyrene		ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Naphthalene	1.1	ug/L		J	1
Q17-W	08-29962-NY02ORE	SW8270D	Phenanthrene	1	ug/L	U	UJ	1
Q17-W	08-29962-NY02ORE	SW8270D	Pyrene	1	ug/L	U	UJ	1
WD-2	08-29963-NY02P	SW6010B	Arsenic		mg/l	U	DNR	11
WD-2	08-29963-NY02P	SW6010B	Lead	0.02	mg/l	U	DNR	11
WD-2	08-29963-NY02P	SW8270D	1-Methylnaphthalene	1	ug/L	U	DNR	10
WD-2	08-29963-NY02P	SW8270D	2-Methylnaphthalene	1	ug/L	U	DNR	10
WD-2	08-29963-NY02P	SW8270D	Acenaphthene	1	ug/L	U	DNR	10
WD-2	08-29963-NY02P	SW8270D	Acenaphthylene		ug/L	U	DNR	10
WD-2	08-29963-NY02P	SW8270D	Anthracene	1	ug/L	U	DNR	10
WD-2	08-29963-NY02P	SW8270D	Benzo(a)anthracene	1.1	ug/L		DNR	10
WD-2	08-29963-NY02P	SW8270D	Benzo(a)pyrene	1.3	ug/L		DNR	10
WD-2	08-29963-NY02P	SW8270D	Benzo(b)fluoranthene		ug/L	U	DNR	10
WD-2	08-29963-NY02P	SW8270D	Benzo(g,h,i)perylene		ug/L	U	DNR	10
WD-2	08-29963-NY02P	SW8270D	Benzo(k)fluoranthene	1	ug/L	U	DNR	10
WD-2	08-29963-NY02P	SW8270D	Chrysene	1.5	ug/L		DNR	10

						Lab		DV Qual
Sample ID	Lab Sample ID	Method	Analyte	Result	Units	Flag	DV Qual	RC
WD-2	08-29963-NY02P	SW8270D	Dibenz(a,h)anthracene	1	ug/L	U	DNR	10
WD-2	08-29963-NY02P	SW8270D	Dibenzofuran	1	ug/L	Ù	DNR	10
WD-2	08-29963-NY02P	SW8270D	Fluoranthene	1.9	ug/L		DNR	10
WD-2	08-29963-NY02P	SW8270D	Fluorene		ug/L	U	DNR	10
WD-2	08-29963-NY02P	SW8270D	Indeno(1,2,3-cd)pyrene	1	ug/L	Ü	DNR	10
WD-2	08-29963-NY02P	SW8270D	Naphthalene	1.8	ug/L		DNR	10
WD-2	08-29963-NY02P	SW8270D	Phenanthrene	2.9	ug/L		DNR	10
WD-2	08-29963-NY02P	SW8270D	Pyrene	2.9	ug/L		DNR	10
WD-2	08-29963-NY02PRE	SW8270D	1-Methylnaphthalene		ug/L		J	1
WD-2	08-29963-NY02PRE	SW8270D	2-Methylnaphthalene	2.2	ug/L		J	1
WD-2	08-29963-NY02PRE	SW8270D	Acenaphthene		ug/L		J	1,9
WD-2	08-29963-NY02PRE	SW8270D	Acenaphthylene		ug/L	U	ÜJ	1
WD-2	08-29963-NY02PRE	SW8270D	Anthracene	1	ug/L	Ü	UJ	1
WD-2	08-29963-NY02PRE	SW8270D	Benzo(a)anthracene	1	ug/L	U	UJ	1
WD-2	08-29963-NY02PRE	SW8270D	Benzo(a)pyrene	1	ug/L	U	UJ	1
WD-2	08-29963-NY02PRE	SW8270D	Benzo(b)fluoranthene	1	ug/L	U	UJ	1
WD-2	08-29963-NY02PRE	SW8270D	Benzo(g,h,i)perylene	1	ug/L	U	UJ	1
WD-2	08-29963-NY02PRE	SW8270D	Benzo(k)fluoranthene	1	ug/L	U	UJ	_ 1
WD-2	08-29963-NY02PRE	SW8270D	Chrysene	1.2	ug/L		J	1
WD-2	08-29963-NY02PRE	SW8270D	Dibenz(a,h)anthracene	1	ug/L	U	UJ	1
WD-2	08-29963-NY02PRE	SW8270D	Dibenzofuran	1	ug/L	Ū	ÚĴ	1
WD-2	08-29963-NY02PRE	SW8270D	Fluoranthene	1.5	ug/L		J	1
WD-2	08-29963-NY02PRE	SW8270D	Fluorene		ug/L	U	UJ	1
WD-2	08-29963-NY02PRE	SW8270D	Indeno(1,2,3-cd)pyrene	1	ug/L	U	UJ	1
WD-2	08-29963-NY02PRE	SW8270D	Naphthalene	12	ug/L		J	1,9
WD-2	08-29963-NY02PRE	SW8270D	Phenanthrene	2.5	ug/L		J	1
WD-2	08-29963-NY02PRE	SW8270D	Pyrene	2.2	ug/L		J	1,9



# APPENDIX C COMMUNICATION RECORDS

From: Chris Ransom

**Sent:** Friday, November 21, 2008 11:02 AM

To: 'markh@arilabs.com'; 'Norm Puri'

Subject: BNSF SDG NM40

#### Hi Mark,

There are two water samples on the COC for NM40, Q2-D-W and WD-1. They were marked for SVOCs, NWTPH-DX, NWTPH-G/BTEX, Diss Metals and turbidity, however we do not have any data for SVOC or NWTPH-Dx. Were these analyses cancelled or perhaps errors on the COC? If you could let me know either way, I'd appreciate it.

#### Thanks

Chris

#### **Christine Ransom**

Sr. Project Chemist EcoChem, Inc. 710 2nd Ave. Suite 660 Seattle, WA 98104 ph 206 233-9332 ext.109 fax 206 233-0114 cransom@ecochem.net State of the standard of

From: Chris Ransom

Sent: Tuesday, November 04, 2008 10:15 AM

To: 'markh@arilabs.com'
Subject: POS BNSF SDG NM43

Hi Mark,

For the metals analyses of the sediment and soil samples in SDG NM43, both matrices were digested in the same prep batch. One of the sediment samples was used for the duplicate and matrix spike. The sediment and soil samples had different analyte lists though, and only the sediment analytes were reported for the method blank, LCS, Dup, and MS. Nickel was a target analyte for two of the soils, but since it was not on the list for the sediments, it was not reported for the QC samples. Could you please resubmit the summary forms for MD, LCS, Dup, and MS with nickel included?

**Thanks** 

Chris

#### **Christine Ransom**

Sr. Project Chemist EcoChem, Inc. 710 2nd Ave. Suite 660 Seattle, WA 98104 ph 206 233-9332 ext.109 fax 206 233-0114 cransom@ecochem.net you walked had

From: Chris Ransom

Sent: Monday, January 05, 2009 4:44 PM

To: 'Norm Puri'

Subject: FW: POS Quendall SDG NY02

Hi Norm,

I just wanted to let you know that these samples are being re-analyzed by ICP-MS in order to achieve the QAPP required reporting limits. We should have the data around the end of the week. Let me know if you have any questions.

Chris

From: Chris Ransom

**Sent:** Monday, December 08, 2008 12:58 PM

To: 'markh@arilabs.com'; 'Norm Puri'; 'Steve Perrigo'

Cc: Linda Bohannon

Subject: POS Quendall SDG NY02

Hi Mark,

We noticed that the water samples in SDG NY02 were analyzed for arsenic and lead by ICP (6010) instead of by ICP-MS (200.8). The results were reported in mg/L (ppm) instead of ug/L (ppb) with much higher reporting limits than are specified in the QAPP. For these Round 2 samples, the results for arsenic are all ND at 0.05 mg/L (50 ug/L) and the results for lead are all ND at 0.02 mg/L (20 ug/L). I believe these samples will need to be reanalyzed by ICP-MS in order to achieve the necessary MRLs of 0.2 ug/L for arsenic and 1.0 ug/L for lead. All Round 1 water samples (SDGs NM40 and NM43) were analyzed by ICP-MS with the appropriate reporting limits.

Could you please let me know as soon as possible if there is sufficient sample remaining for re-analysis and what the TAT would be?

Thank you,

Chris

#### **Christine Ransom**

Sr. Project Chemist EcoChem, Inc. 710 2nd Ave. Suite 660 Seattle, WA 98104 ph 206 233-9332 ext.109 fax 206 233-0114 cransom@ecochem.net

From:

Chris Ransom

Sent:

Wednesday, December 17, 2008 12:17 PM

To:

'markh@arilabs.com'

Cc:

'Sue Dunnihoo'; Linda Bohannon

Subject: POS BNSF Diesel data

Tracking: Recipient

Delivery

'markh@arilabs.com'

'Sue Dunnihoo'

Linda Bohannon

Delivered: 12/17/2008 12:17 PM

#### Hi Mark,

We seem to have found a problem with the diesel calculations in some of the POS BNSF data packages. The same curve affects SDGs NX66, NX71, NX79, and NY02. The diesel calibration curve was analyzed on 11/4/08 on instrument FID3A.I. The average RF value for the WA diesel range is 17742 in the summary table, however the samples were calculated using an RF value of 16911.5, which appears on all of the quant reports. The motor oil and surrogates all use the correct RRF values from this ICAL. Could you check into this and see if perhaps the WA diesel RRF value wasn't updated based on this ICAL?

Also, have you had a chance to check into the question regarding the metals data for SDG NY02? The water samples were analyzed for arsenic and lead by ICP and have much higher detection limits than specified in the QAPP. The samples may need to be re-analyzed by ICP-MS in order to achieve the necessary detection limits.

Please feel free to give me a call if you need any more information.

Thanks

Chris

Christine Ransom

Sr. Project Chemist EcoChem, Inc. 710 2nd Ave. Suite 660 Seattle, WA 98104 ph 206 233-9332 ext.109 fax 206 233-0114 cransom@ecochem.net nuve submitted

From:

Sue Dunnihoo [sue@arilabs.com]

Sent: To: Friday, January 02, 2009 5:27 PM Chris Ransom; Mark Harris

Subject:

**POS BNSF** 

Attachments:

TPH\_Curve\_102908.pdf



TPH\_Curve\_102908 .pdf (780 KB)

Here's a copy of the curve that all the data was evaluated against. I think this should solve the issue without having to reprocess all the data against the other curve and redo all the reports, since the CCALs passed and were evaluated against this curve.

- Sue

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